# **tyFLOW** Help Quick Jump Page

## CLICK ON OPERATOR TO QUICK JUMP TO THAT HELP PAGE

	Birth	8	Object Bind	8	Link to Target	R	Cloth Bind		Send Out
0,0	Birth Burst	380	Particle Force	0+0	Move to Target		Cloth Collect	2	Split
	Birth Flow	00	Particle Groups	0	Set Target		Modify Bindings	<u> </u>	Surface Test
25	Birth Fluid	3	Path Follow	10	Instance ID	100	Particle Bind	(E)	Time Test
	Birth Intersections	0	Point Force		Instance Material	(8)	Particle Break	558	Camera Cull
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×	Delete		Speed	4	Tets		Actor Animation	8	VDB Solver
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HELP FILE DATE: May 19 2020

### **HELPERS QUICK JUMP LINKS**



### **MODIFIERS QUICK JUMP LINKS**

tyUVWFreeze	tyCarve	tyConform	tyEdgeWeights	tyFaceExtrude	tyFaceFracture	tySmooth	tySelect
tySplineMesher	tyWeld	tyBoolean	tySplineExtrude	tyParticleSkin	tyProperties	tySpiral	tySlice
tyBend			tySplineCache				

#### **MISCLEANEOUS LINKS**

MAXScript	tySpaceWarp	tyFlow Controllers	tySplinesObject
tyPreviewUtility	tyMaterial	Texmaps	

## **Understanding the Simulation Loop**

**tyFlow's** simulation loop is evaluated in the following way:

For each time step of the simulation:

- 1) The optimal order of events in the flow is found, by doing a depth-first search of connected events, starting with events that contain birth operators.
- 2) The ordered events are processed in ascending order, by evaluating the event's operators from top to bottom.

**NOTE:** Certain operators may have their evaluation deferred until later in the loop. Particle Physics operators set to "integration" mode are only evaluated during the bind solver step. Collision operators are also only evaluated after all other operators and non-PhysX solvers have been evaluated.

- 3) Once the events have been processed, the bind solver is evaluated. Cloth binds, particle binds, and any Particle Physics operators set to "integration" mode are evaluated.
- **4**) Once the bind solver is finished, the wobble solver is evaluated. The wobble solver calculates spring forces applied to particles created with a Wobble operator.
- 5) Once the wobble solver is finished, the actor solver animates the transforms of any actor rig particles that have animation clips applied to them.
- **6**) Once the actor solver is finished, all accumulated particle velocity and spin values are integrated into the system.
- 7) After velocity/spin integration, PhysX objects are processed. Internal PhysX rigidbodies have their position/velocities matched to their corresponding particles and the PhysX solver is evaluated.
- 8) Once the PhysX solver is finished, some final adjustments are made to the simulation. The bind solver checks to see if any particles should be put to sleep, the age of all particles is incremented, appropriate particles are cached, etc.

## **Simulation Validity**

Under normal circumstances with moving particles, the simulation is history-dependent. This means that in order for the simulation state at time [t] to be known, the simulation state at time [t - timestep] must also be known. This also means that any given simulation state in those circumstances will only be valid for one timestep. As soon as the time changes, the simulation state changes, rendering the previous state invalid.

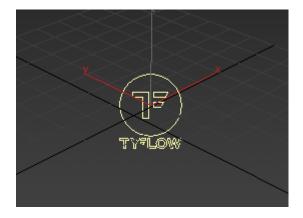
The exception to this case is if there are no changing properties in the simulation. If nothing is changing, then the simulation state at all times will be the same because the simulation is static.

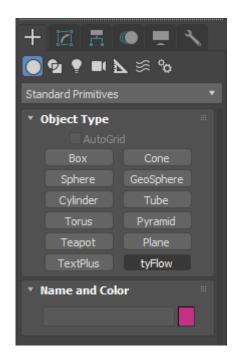
When the simulation is initialized, **tyFlow** iterates over all operators and queries them for the types of changes they will make to the simulation. If every single operator reports that it does not make changes to particles over time, then **tyFlow** will recognize that the simulation will remain static forever once all new particles have been birthed. This makes **tyFlow** a very efficient geometry scattering tool, since static particles that are scattered in a scene will not have to be continually evaluated after they are birthed.

**NOTE:** Grey text labeled: "Static: [range]" appears in the bottom left corner of the editor and displays info about which frames of the simulation are static, or tells you if the simulation is never static.

# Creating a new tyFlow particle system

To create a **tyFlow** scene object, navigate to Create—Standard Primitives—**tyFlow** within the modifier panel and choose a location in the scene to place the object.





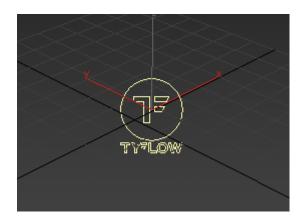
**NOTE:** The location you choose to place your **tyFlow** object does not matter – all **tyFlow** calculations happen in World-Space (not Object-Space), so the object's transform will have no effect on the simulation or display of the particles.

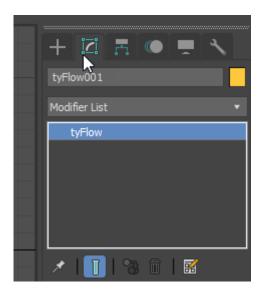
## **Using the Editor**

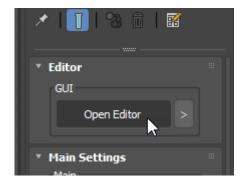
The editor is where flows reside within a **tyFlow** object.

## Opening the Editor

With the tyFlow icon selected in any viewport, go to the modifier panel and scroll down to open editor and click the button







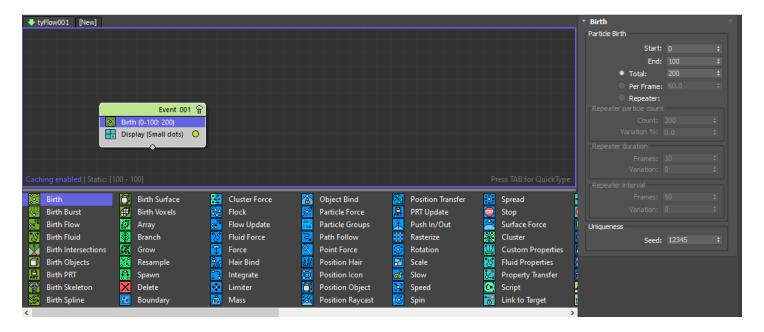
To open the editor for a particular **tyFlow** object, select 'Open Editor' from the modifier panel or "Editor..." for a particular object in the **tyFlow** viewport menu.

**NOTE:** Unlike *Particle Flow*, which features a single editor for all flows in the scene, **tyFlow** objects each have their own editor.

**NOTE:** The size and position of each editor window is saved to the scene file, and will be restored when loading the scene. If an editor is saved to an off-screen location, it will be automatically centered inside the current window

## **Navigating the Editor**

By using your middle mouse button and scroll wheel, you can pan and zoom in the editor. Left-click can be used to select events/operators/connections/etc, and right-click can be used to display relevant context menus.



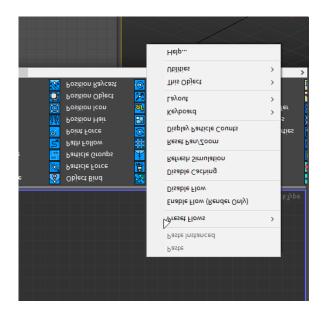
The grid view is where all flows are constructed. The operator list below the grid view displays all available operators. The rollout panel to the right of the grid view displays all settings for any operator that is selected.

NOTE: The pan and zoom value of each editor window is saved to the scene file, and will be restored when loading the scene.

## Right-click Menus

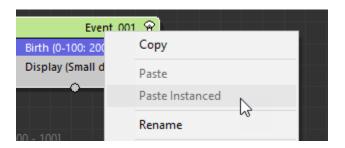
Various right-click context menus will appear when right-clicking the grid view, events, or operators.

While most right-click menu options are self-explanatory, a few may require some further details:



### Operator right-click menus:

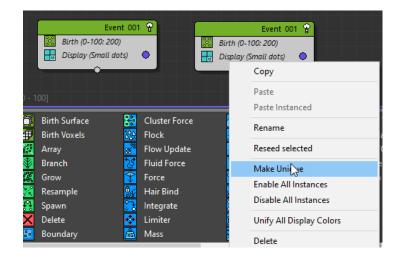
• **Paste Instanced**: operators that are pasted as instances will share all settings with the operator they were copied from. This menu item will only appear if an operator has previously been copied.



• Make Unique: choosing this option on an operator that is an instance of another operator will make its settings fully independent from that other operator.

Therefore, from that point on it will no longer be an instance of the other operator.

This menu item will only appear when right-clicking on an instanced operator.



• Make Instances: if multiple operators are selected, the operator that was right-clicked will become the data source of all the other operators which are selected. In other words, all operators that are selected will lose their current settings and be made instances of the operator for which this option is chosen. This menu item will only appear if multiple operators of the same type are selected.

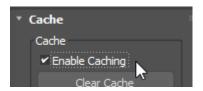
### TIP

"Make instances" is an easy way to transfer settings between existing operators. For example, if you have two independent operators of the same type and you want them to have the same settings, instead of deleting one and copy/pasting the other in its place, simply select them both and choose "make instances" on the one whose data you want to copy to the other. The other operator will be instantly converted into an instance of the one you right-clicked. In that sense, "make instances" is shorthand for "make all of the other selected operators of the same type instances of this one".

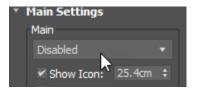
#### **Editor hints**

The editor displays various bits of information around its frame to assist users in diagnosing different behaviors of the flow.

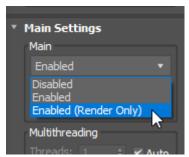
A **blue outline** around the frame of the editor means that realtime caching is enabled. "Caching enabled" will also appear in the bottom left corner.



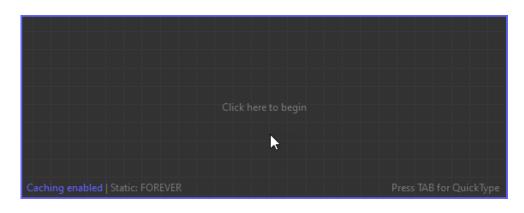
A **red outline** around the frame of the editor means that the flow has been disabled. "Simulation disabled" will also appear in the bottom left corner.

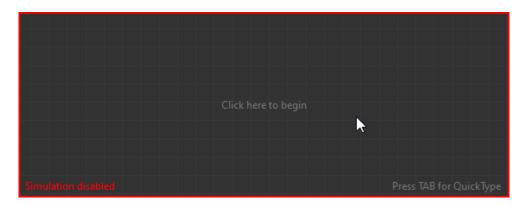


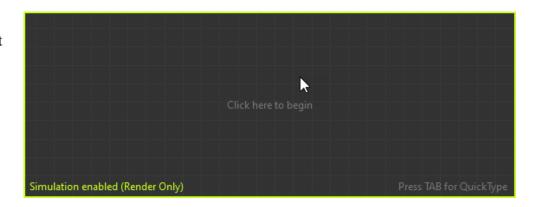
A **yellow outline** around the frame of the editor means that the flow is in render-only mode. "Simulation enabled (Render Only)" will also appear in the bottom left corner.

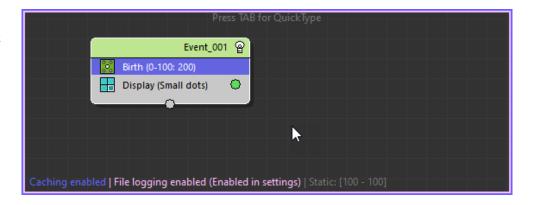


A **pink outline** around the frame of the editor means that the flow is logging its simulation progress to disk. Relevant information about the location of the log will also appear in the bottom left corner.



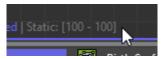








Grey text labeled "Static: [range]" appearing in the bottom left corner displays info about which frames of the simulation are static.



**NOTE:** A static frame is a frame which requires no additional simulation steps because it contains particles which do not change over time. A completely static flow with no moving particles will only need to evaluate a single simulation step, allowing for fast updates and timeline scrubbing.

## **Operator Category Tabs**

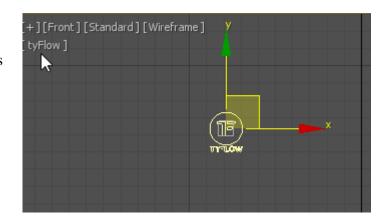


Below the grid window is a tab selection window, you can click on the different operator color groups to filter the list of operators shown in the operator window.

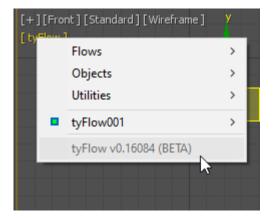
This allows you to quickly browse color coded categories.

## Viewport Menu

tyFlow features a viewport menu (located in the top left corner of the active viewport) that will automatically be displayed when any tyFlow objects are present in the scene, and will be automatically hidden when no tyFlow objects are present in the scene. It can be used to quickly access and control available tyFlow objects, with options to enable/disable, hide/unhide, refresh, select and edit them. Options also exist to select/hide/unhide all existing tyFlow objects at once.



The viewport menu also displays the current version number of the **tyFlow** plugin file installed on the machine.



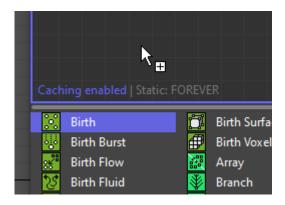
**NOTE:** Using the viewport menu to access **tyFlows** in the scene is often much easier than manually navigating to them through the viewport or scene explorer, and is part of an efficient workflow.

## **Creating Flows**

Flows are groupings of operators, events and connections between them. They allow you to direct the behavior of particles over time.

#### **Creating Operators**

To create a new operator, drag an operator from the operator list into the grid view.



An operator dragged into an existing event will be added to that event's operator list.



An operator dragged directly over the grid will be added to a new event. Operators can also be copied and pasted (from their right-click menu), and the same rules apply – operators pasted over the grid will be assigned to a new event, and operators pasted into an existing event will be added to that event's operator list.

**NOTE:** You can copy and paste multiple events and/or operators at the same time, by selecting them with the drag marquee, or by holding CTRL to individually select more than one at a time.

#### **Creating Events**

Events themselves cannot be created directly – only copied and pasted (from their right-click menu) or created by a dragging an operator over the grid.

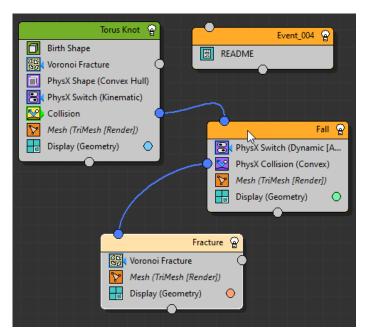
**NOTE:** When a new event is created it will automatically be assigned a new *Display* operator, if necessary.



#### **Creating Connections**

Connections between operators and events allow you to direct the behavior of particles over time. If an operator's output is connected to an event's input, any particle that satisfies the test condition of the operator will be sent to the connecting event at the end of the operator's simulation step. The direction of a flow is always forward (from operator to event) – operators can send particles to events, but events cannot send particles back into operators.

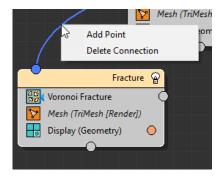
An event can take inputs from multiple operators, but an operator cannot output particles to multiple events.



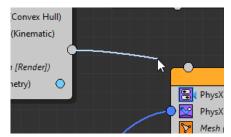
**NOTE: tyFlow** correctly handles event looping, where an operator is connected to a prior event, such that particles loop back to that same operator within the same timestep (normally resulting in an infinite loop that can never complete). No extra measures need to be taken to avoid infinite looping in those cases, as **tyFlow** will automatically ensure only a single loop is completed per timestep in such a scenario

#### **Shaping connection wires**

By right-clicking on a connection wire, you can add or remove points to it.



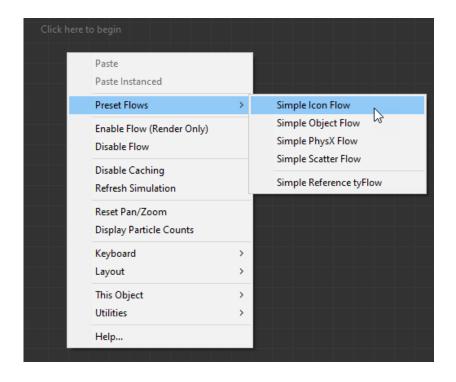
By dragging those points, you can shape the wire, allowing you to visually route wires around events in the grid.



#### **Preset Flows**

In the right-click context menu of the grid view, you can find a "New" submenu that lists several preset flows which can be created.

Creating a new preset flow will *not* reset the editor or affect existing flows – it will merely *add* the selected preset flow to the editor. Relevant scene objects used to control preset flow properties will also be created, depending on the preset chosen.



# tyFlow Object Settings

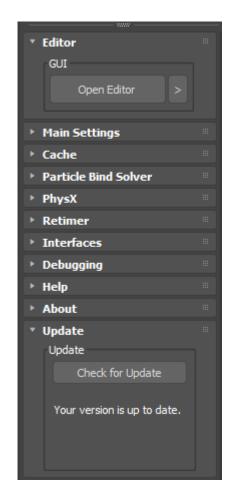


**tyFlow** objects are what contain the events and operators requires to create and compute particle simulations.

Each **tyFlow** object contains a variety of global settings which can be used to tune simulations, or export particles to various formats.

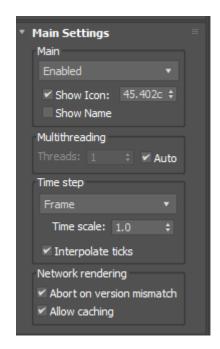
These settings can be accessed within the modifier panel.

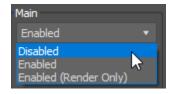




We will go through each rollout category in the sidebar to find information about a particular setting.

## **Main Settings Rollout**





**Enabled state dropdown**: controls whether the flow is enabled, enabled (Render Only), or disabled. Disabled flows do not evaluate at all, while enabled (Render Only) flows only evaluate at render time.

Show icon: controls whether the **tyFlow** object's icon is visible in the viewport.

Icon size: controls the size of the **tyFlow** object's icon in the viewport.

Show name: controls whether the name of the **tyFlow** object is displayed in the viewport.

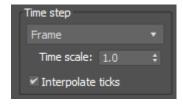
## Multithreading

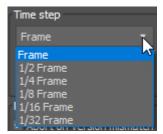
Thread count: controls the maximum number of CPU threads the flow can use to evaluate the simulation.

Auto: allows **tyFlow** to determine the maximum number of threads to use to evaluate the simulation (defaults to max. available)

**Note:** Setting thread count to a particular value doesn't mean that number of threads will be used for every operation, only that **tyFlow** may not exceed that particular number of threads for a given operation. Some operations benefit from more threads and some with less, and **tyFlow** makes internal determinations regarding the actual number of threads to use on a per-algorithm basis (with the only constraint being the maximum value provided by the user here).

#### Time Step





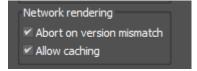
Time step dropdown: controls the number of steps the simulation will divide each frame into. With more time steps, simulation accuracy will be increased but simulation speed will be decreased. A value of 'Frame' is usually fine for simulations that don't require physical accuracy. A value of "1/4 Frame" or "1/8 Frame" is best for simulations featuring physically accurate constraints (sand/cloth/rope/etc) in order to increase overall simulation stability.

Time scale: 1.0 † Time scale: the time scale setting allows you to control the speed of the simulation. Unlike the retimer settings, which allow you to control playback speed *after* the simulation is cached, the time scale setting allows you to control simulation speed *as it is calculated*. Values less than one have the effect of slowing the simulation, and values greater than one speed things up.

**Note:** Increasing the time scale value can lead to physics/PhysX inaccuracies, because increasing the value increases the velocity of all particles within the time step. You would only use this setting, as opposed to using the retimer, if you need to change the simulation speed of particles while maintaining the speed of all scene objects your particles are interacting with.

Interpolate ticks: controls whether particle transforms will be interpolated at ticks between time steps.

#### **Network rendering**



**Abort on version mismatch**: Rendering **tyFlows** on machines whose **tyFlow** version does not match the version of the originating scene file is usually a bad idea. While it doesn't guarantee problems, if the scene file relies on a feature not present in the version on the render machine, the scene may not render correctly. If this setting is enabled, renders will be automatically aborted on machines whose **tyFlow** version does not match the scene file, avoiding incorrect renders in the process.

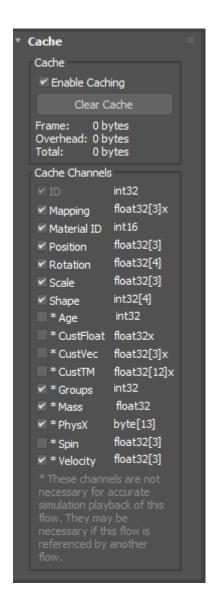
Note: It is always recommended to use the most recent version of tyFlow.

**Allow caching**: Controls whether render machines will cache frames while the simulation is processed on them prior to rendering. In most cases this should be left on, and only disabled if a particular render machine has enough RAM to process the simulation, but not enough RAM to cache each processed frame. Disabling this can greatly increase render time of complex simulations if multiple frames of the simulation are accessed in descending order (for example, in the case of a simulation retimed to play in reverse, or multiple flows referencing each other with staggered frame offsets).

## **Cache Settings Rollout**

**tyFlow's** realtime timeline caching allows for smooth playback of a simulation in the viewport after it has been computed.

**Note:** Caches are saved on a per-frame basis, and do not adhere to the simulation's time step setting (the simulation itself will still run at whatever time step interval is chosen, but the cache will only save values once per frame). **tyFlow** will still interpolate the subframes of cache data if the flow's "interpolate ticks" setting is enabled, but if you need to playback your flow with sub-frame accuracy, you should disable caching.



### Cache

**Enable Caching**: controls whether realtime timeline caching is enabled or disabled.

**Clear Cache**: clears the contents of the cache and effectively resets the simulation.

#### Cache channels

[Channel name - data type]: lists the available channels to save with the cache, and their corresponding data type.

Data types and their corresponding size in bytes:

byte = 1 byte

int16 = 2 bytes

int32 = 4 bytes

float16 = 2 bytes

float32 = 4 bytes

**Note:** The number in square brackets next to some data types represents the number of values that must be stored for that particular channel. For example, a position channel requires X/Y/Z values, each of which is stored as a float32 data type. So the size in bytes for a particular particle position value is 12 bytes (float32 x 3 values).

Mapping values are stored as a float32[3] x the number of mapping channels assigned to the particle.

Custom float data values are stored as a float32 x the number of custom float data channels assigned to the particle.

Custom vector data values are stored as a float32[3] x the number of custom vector data channels assigned to the particle.

Custom TM data values are stored as a float32[12] x the number of custom TM data channels assigned to the particle.

#### TIP:

You can estimate how much RAM will be required to cache a particular flow by using the following equation:

### (number of particles) x (number of frames) x (total size of all saved channels in bytes)

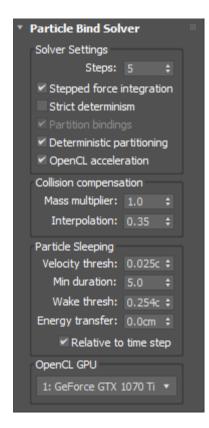
For example, a simulation of 1 million particles over 250 frames which caches particle positions and velocity will require approximately 7GB of RAM. This resulting value should only be considered an estimate, since caching requires some extra overhead not contained within any particular channel. The amount of overhead varies depending on the complexity of the flow and is displayed in the Cache rollout.

#### **WARNING:**

Caching can use a lot of RAM up very quickly, depending on the complexity of your flow. If your machine has limited RAM and you are planning on simulating many millions of particles, it is best to turn caching off and instead export the particles to disk using **tyFlow's** available export options. **tyFlow** makes **no** effort to check whether RAM allocations are possible on a given system, which means that if your RAM is full and **tyFlow** needs more of it in order to continue a simulation, **tyFlow** could cause crash. This 'fast and loose' approach to simulating is by design, and it is up to the user to ensure their system is capable of handling the simulations they are trying to run.

## **Particle Bind Solver Settings Rollout**

**tyFlow's** particle bind solver is what solves all inter-particle bindings (aka **constraints** or **joints**) within **tyFlow** (excluding PhysX bindings). At its core, a binding is just a relationship between two particles, and solving many bindings in succession is what gives rise to intricate behaviors seen in materials like dirt, wet sand, cloth, ropes, etc. Proper tuning of the bind solver is important when simulating these complex systems.



## **Solver Settings**

Steps: 5 \$ Steps: controls the number of sub steps per simulation time step to solve all active bindings.

**Note:** The total number of evaluations per binding per frame can be calculated as (bind steps) x (simulation time steps). The higher the total number of evaluations, the more accurate the solver results will be. For granular simulations, a simulation time step of either "1/4 Frame" or "1/8 Frame" with bind solver steps of 5-10 is often adequate. For hires cloth simulations, bind solver steps may need to be much higher in order to maintain cloth stiffness.

Stepped force integration: controls whether particle velocities are smoothly added to the bind solver per step, or added only once prior to all bind solver steps. Keeping this enabled has a very minor performance impact but can reduce high velocity artifacts in the resulting simulation.

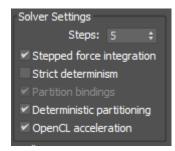
Strict determinism: with this setting turned on, successive runs of a simulation should return identical results. With this setting off, there is no guarantee that bindings will be evaluated in the same order or that race conditions between multiple threads will be prevented, and so results across multiple simulations may vary. Turning this setting on can have a detrimental performance impact, so it's recommended to keep it off, unless you need the simulation to produce identical results across successive runs.

**WARNING:** If you plan on rendering across multiple computers, "strict determinism" must be enabled or else the frames returned by different machines will not be in sync. An alternative to rendering with determinism on is to instead cache out your particles locally and then render a tyCache/PRT/etc loader instead of the **tyFlow** object itself.

If you choose to render your **tyFlow** with "strict determinism" on across multiple machines, make sure all machines have consistent OpenCL support. If you have OpenCL acceleration enabled but not all machines that you are using support OpenCL, mixing CPU/GPU solvers can impact determinism *even with "strict determinism" enabled.* 

#### TIP:

It is generally a better practice to render a cache of your flow instead of your **tyFlow** object itself, when rendering across multiple machines. Rendering a cache ensures that hardware differences between computers have no impact on the consistency of the final output.



## Solver Settings (continued)

**Partition bindings**: controls whether bindings will be split into non-overlapping groups, before being solved in a multithreaded manner. Turning this setting off can decrease simulation time, at the cost of increased error accumulation over time. This setting has no effect when OpenCL acceleration is enabled, because OpenCL acceleration requires partitions.

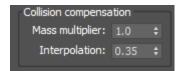
**Deterministic partitioning**: controls whether the partitioning process must avoid threaded race conditions. This setting can usually be disabled for granular flows, but should usually be enabled for cloth/soft-body flows to avoid jittering artifacts.

**OpenCL acceleration**: if an OpenCL2.0-compatible GPU device is found on the system, this option will be available. OpenCL acceleration can increase simulation performance, depending on the power of the available GPU. When enabled, all bindings will be solved on the GPU instead of the CPU.

**Note:** Enabling OpenCL acceleration does not guarantee a performance boost. There is a fair amount of overhead involved in transferring data to-and-from the GPU during the simulation, that can offset the actual speed boost the GPU offers during its calculation phase. While an overall increase in performance should be expected for very high-end GPUs on systems with few CPU cores, a system with many CPU cores and a lowend GPU may not see much of a performance boost with OpenCL at all. Results will vary across hardware and users must experiment to determine if OpenCL acceleration is right for them. It is not a magic bullet solution.

## **Collision Compensation**

During each simulation step, collisions are always processed after bindings. No solid geometry collisions are processed while the bind solver evaluates bindings each bind solver step. Because of this, it's possible for the bind solver to pull particles straight through colliders, only for the subsequent collision step to fix those intersections afterwards. However, even though those intersections are eventually fixed, the rest of the bindings remain unaware that such a collision ever took place, and this can cause visual artifacts within the overall bind network. To compensate for this, particle masses can be artificially adjusted when collisions are detected on the previous simulation substep. Collided particles can be given a heavier mass, so that they won't be pulled as forcefully by their connected particles. Once previously-collided particles are determined to have no more collisions, their mass will return to normal. The combination of these effects can help reduce visual artifacts in binding networks (like cloth).

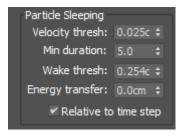


**Mass multiplier**: The multiplier applied to the (inverse) mass of particles that collided on the previous simulation step. The smaller the value, the less influence surrounding bindings will have on a collided particle.

**Interpolation**: The interpolation speed used to transition particle masses between the collision compensation value and their original value. Keeping this value low can help prevent jittering artifacts caused by the masses of collided particles switching between the compensation value and their original value too quickly.

## **Particle Sleeping**

By enabling particle sleeping, you can force low-velocity particles to come to a standstill when they would otherwise keep moving over time. This can prevent unwanted motion in particles and forcibly bring jittering particles to rest.



**Velocity thresh**: particles whose velocity magnitude is below this threshold will be considered candidates for sleep.

**Min duration:** candidate particles whose velocity magnitude remains under the velocity threshold for this duration of time will be put to sleep.

**Wake thresh**: sleeping particles whose velocity exceeds this value at the end of a time step will be awoken.

**Energy transfer**: the amount of neighbor-particle energy that can contribute to waking a particle.

**Relative to time step**: Multiplies threshold velocities by the time step.

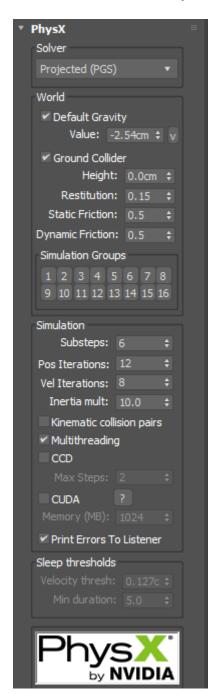
#### INFO:

Because velocities are integrated each time step, wake/sleep thresholds may be too large by default if your time step is less than 1. For example, if your gravity strength is -1.0 and your sleep threshold is 0.5, particles will not fall asleep when your time step is 1 frame. However, if your time step is  $\frac{1}{2}$  frame, particles will fall asleep because at each substep their velocity is increased by 0.5 instead of 1.0 (which matches the sleep threshold). If "relative to time step" is enabled, the wake/sleep thresholds will be multiplied by the time step delta, and so in this example the effective threshold would actually be 0.25 (0.5 \*  $\frac{1}{2}$ ) per step.

**NOTE:** Particle sleeping has no performance impact. Its impact is purely visual. Sleeping particles will still be evaluated by the solver – the difference is that if they are considered asleep at the end of a time step, they will be returned to their previous location (effectively rendering them motionless).

## **PhysX Rollout**

These controls affect all PhysX rigid bodies in the simulation.



#### Solver

**PGS/TGS**: controls which PhysX solver to use in the simulation

**INFO:** Information about each solver (Projected/Temporal Gauss-Seidel) can be found on NVidia's PhysX website. TGS is a relatively new addition to PhysX, and is typically faster than PGS for rigid body simulations, but can produce unexpected results when solving PhysX bindings (constraints). In general, it should be treated as an experimental solver, and PGS should still be used by default.

#### World

**Default gravity**: controls whether a default gravity force will be applied to all PhysX particles.

Gravity value: controls the strength of the default gravity force.

**Ground collider**: controls whether a default ground collider will be added to the PhysX simulation.

**Height**: the height of the default ground collider, in world-space.

**Restitution**: the restitution of the default ground collider.

**Static friction**: the static friction of the default ground collider.

**Dynamic friction**: the dynamic friction of the default ground collider.

**Simulation groups**: the simulation groups that will be affected by the ground collider

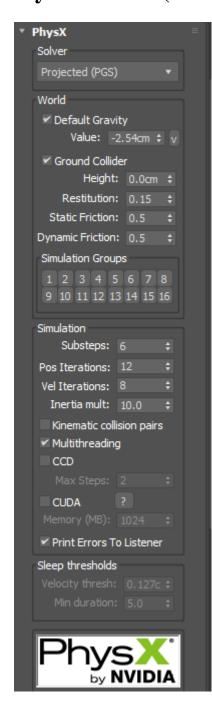
#### Simulation

**Substeps**: the number of substeps that will be calculated per time-step for the PhysX simulation.

#### TIP:

Increase substeps to increase the overall accuracy of the simulation. For high-fidelity simulations, a value of 12 or higher may be more appropriate than the default value.

## **PhysX Rollout (continued)**



### Simulation (continued)

**Pos iterations**: the number of position iterations that will be used to solve joint/contact constraints, per rigidbody, per substep.

**Vel iterations**: the number of velocity iterations that will be used to solve joint/contact constraints, per rigidbody, per substep.

#### TIP:

Increase pos/vel iterations to reduce jittering when simulating particles with PhysX bindings.

**Inertia mult**: this is a multiplier which affects all PhysX particles' inertia. Higher values can increase simulation stability while decreasing angular acceleration/deceleration. Lower values increase angular acceleration/deceleration, but can lead to jittering or other instabilities. For smaller objects with low mass, this value can be lowered (0.5 - 1.0). For bigger objects with a lot of mass, it should be kept high (5.0 - 20.0).

**Kinematic collision pairs**: controls whether inter-penetrating particle rigidbody pairs that are set to 'kinematic' or 'trigger' will generate contacts.

**Multithreading**: controls whether the PhysX engine will make use of multiple CPU threads.

**CCD**: controls whether *continuous collision detection* is enabled or disable. *Continuous collision detection* can prevent collision tunnelling between high-velocity rigidbodies, for a performance cost.

**CCD steps**: controls the number of substeps used to resolve collisions by the CCD engine.

**CUDA**: controls whether PhysX computations will be accelerated with CUDA.

#### **INFO:**

In order for CUDA acceleration to work, tyFlow requires that two DLL files (PhysXDevice64.DLL and PhysXGPU64.DLL - both available on the tyFlow download page) be placed in the same folder where the tyFlow DLO file is loaded from.

**NOTE:** CUDA acceleration does not always guarantee faster simulations. Due to performance costs related to transferring necessary data to-and-from the GPU, speed benefits from CUDA might not be apparent until hundreds/thousands of rigid body particles are in the simulation. When a simulation has only a few rigid body particles, CUDA acceleration being enabled may actually decrease overall performance.

#### **INFO:**

Setting the CUDA memory limit higher than the default value does not mean the simulation will run faster. The memory limit controls the amount of VRAM to allocate for constraint/contact processing, and generally a CUDA simulation does not require much VRAM in order to process all contacts, even if a lot of rigid bodies are present in the simulation. Setting this value very high is usually unnecessary, and can actually contribute to slowdowns at the beginning of the simulation, due to the time it takes to initially allocate the VRAM. Just because you have a GPU with a lot of VRAM, does not necessarily mean you should increase this setting from its default value. For reference, the default value suggested by NVidia is approximately 140mb.

Print Errors To Listener

Print errors to listener: controls whether to print PhysX simulation errors (reported internally by the PhysX engine) to the MAXScript listener.

### Sleep thresholds



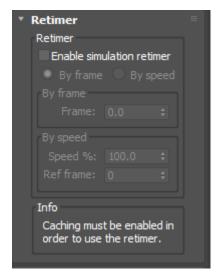
**Velocity thresh**: Particles with a linear/angular velocity below this threshold at the end of the time step will be candidates for sleeping.

**Min duration**: Particles which satisfy the velocity threshold for this number of frames will be put to sleep.

**NOTE:** Strange behavior can occur if particles with PhysX Bindings are put to sleep, therefore particles with PhysX Bindings will ignore the sleep threshold settings.

## **Retimer Rollout**

The **tyFlow** retimer can be used to retime an entire simulation. By animating the retimer playback frame value, you can choose what part of the simulation plays back over the course of the timeline. Subframe values will be correctly interpolated by the retimer, allowing for smooth slow-down or speed-up effects.



**Enable simulation retimer**: controls whether simulation playback will be controlled by the retimer frame value.

**Retime type**: controls whether the retimer affects playback frame or speed.

### By frame

**Frame**: the retimer frame value that controls simulation playback. Animate this value to control the current playback frame.

### By speed

**Speed %:** the percent value that controls simulation playback speed.

**Ref Frame**: the reference frame that the speed multiplier will be relative to.

#### **INFO:**

Setting a proper reference frame is important for the speed multiplier. The reference frame should typically be the start frame of your playback sequence, not necessarily the start frame of the simulation.

The equation for the speed retimer is:

playbackFrame = [referenceFrame + abs(time - referenceFrame) \* speed \* sign(time referenceFrame)].

**TIP:** The speed value is not animatable. If you need a variable playback speed, use the retimer in "frame" mode instead.

## **Interfaces Rollout**

Each **tyFlow** is able to export "I\_PARTICLEOBJ" and "I\_VRAYGEOMETRY" interfaces to 3ds Max. In some cases, enabling these can cause issues between **tyFlow** and other plugins. In other cases, you might want to limit which particles are sent back to a particle query request. These settings allow you to override these interface exports.



## ParticleObject Interface

**Enable Particle Interface**: controls whether the particle interface will be returned when the **tyFlow** object is queried for its default interface. Returning this interface allows third-party plugins to query a **tyFlow** for its particles.

**Incl. legacy interface**: when enabled, 3ds Max's legacy particle interface (I\_PARTICLEOBJ) will be enabled, as well as its modern interface (IParticleObjectExt).

**Note:** Disabling the "incl. legacy interface" setting can improve compatibility with certain 3rd party plugins.

**Export groups**: controls which particle export groups will be returned when a **tyFlow** is queried for its particles. Use these groups to limit which particles will be sent to third-party plugins querying a **tyFlow** for its I PARTICLEOBJ interface.

#### TIP:

For example, if you have a PhoenixFD grid which uses a **tyFlow** as a PHXSource input object, you might only want to use certain particles in the flow to affect the PhoenixFD simulation. By limiting the particles exported through the I\_PARTICLEOBJ interface (by setting the desired export groups), you can limit which particles will be returned to the PhoenixFD object.

## Particle scale queries

When a 3rd party plugin asks tyFlow for the scale of a particle, these options let you choose what property of the particle is returned.

**Return scale**: the explicit scale value of the particle will be returned.

**Return shape diameter**: the diameter of the particle's shape mesh will be returned.

### **VRay Interface**

**Enable VRay Interface**: controls whether the I\_VRAYGEOMETRY interface will be returned when the **tyFlow** object is queried by VRay for its VRay-compatible interface.

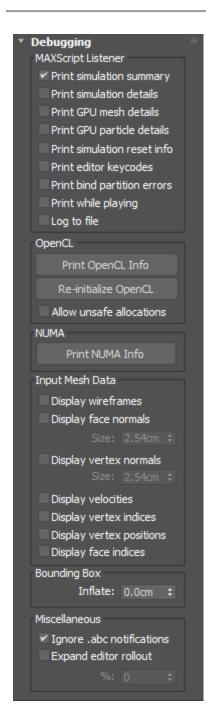
# **Help Rollout**

The Help rollout contains links to the official **tyFlow** documentation, as well as the official **tyFlow** discussion forums.



## **Debugging Rollout**

This rollout contains controls which allow users to profile and debug various aspects of a flow.



### **MAXScript Listener**

**Print simulation summary**: controls whether a summary of flow details will be printed to the MAXScript listener each time a sequence of frames is simulated.

**Print simulation details**: controls whether a verbose list of simulation timing details will be printed to the MAXScript listener each time a sequence of frames is simulated.

**Print GPU mesh details**: controls whether a verbose list of timing details will be printed to the MAXScript listener regarding the time required to upload **tyFlow** meshes to the GPU for display.

**Print GPU particle details**: controls whether a verbose list of timing details will be printed to the MAXScript listener regarding the time required to upload **tyFlow** particle transforms to the GPU for display.

**Note:** Simulation details can provide useful insights into the time required to complete each step of the simulation process. If a simulation is running slowly, you can view the simulation details to see which operator or function is taking the most time to process.

Computing simulation details has a minor performance cost, and printing simulation details to the listener is a slow process, so these settings should remain disabled unless you are actively trying to debug or profile a simulation.

**Print simulation reset info**: controls whether information about changes to input objects is printed to the MAXScript listener.

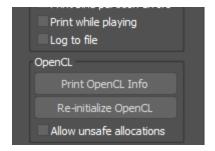
#### TIP:

Every time a flow's input objects change, the flow's simulation is reset to account for the changes. Sometimes a buggy input object may send rogue notifications to the flow, announcing it has changed, even though hit has not. This can cause a flow to continually reset its simulation in an undesirable manner. Enabling this setting can help users figure out which objects are sending change notifications to the flow.

**Print editor keycode**: keycodes of keys pressed in the editor will be printed in the MAXScript listener.

**Print bind partition errors**: if a bind partition contains adjacent constraints, an error will be printed to the MAXScript listener.

**Note:** Partition error printing is a developer setting that should generally be kept off - it involves extra calculations that will slow down the simulation, and doesn't provide any useful information to regular users.



**Print while playing**: controls whether simulation summaries or details will be printed to the MAXScript listener during timeline playback. Keeping this setting disabled will maximize viewport playback speed, by suppressing messages while playback is occurring.

**Log to file**: simulation profiler data will be saved to a log file on the hard disk.

### OpenCL

**Print OpenCL Info**: pressing this button will print a list of available OpenCL devices and their properties to the MAXScript listener.

**Re-initialize OpenCL**: rebuilds OpenCL programs and kernels. In case of an OpenCL error during simulation, OpenCL must be re-initialized before it can be utilized again.

**Allow unsafe allocations**: allows OpenCL to try to allocate more VRAM than the default allocation limit (which is ½ of total VRAM), where necessary.

## **Input Mesh Data**

These settings allow users to see visual properties of a flow's input meshes in the viewport. The resulting markers represent the exact data held in RAM by the flow in its custom **tyMesh** object format, not necessarily the raw data contained within the input objects' Mesh objects. Usually, there should not be a discrepancy between the two, but these options will allow you to see if there is.

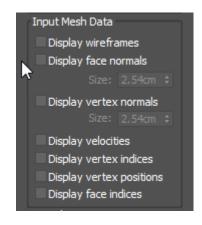
**Display wireframes**: manually draws the wireframes of all input meshes in the viewport.

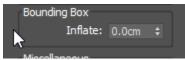
**Display face normals**: manually draws the face normals of all input meshes in the viewport.

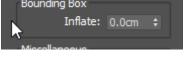
**Face normal size**: controls the overall length of the drawn normals.

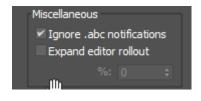
**Display velocities**: manually draws vertex velocity vectors of all input meshes in the viewport.

**Display vertex indices**: manually draws vertex index numbers for all faces of all input meshes in the viewport.









**Display vertex positions**: manually draws vertex position values of all input meshes in the viewport.

Display face indices: manually draws face index numbers for all faces of all input meshes in the viewport.

## **Bounding box**

**Inflate**: manually inflates the bounding box of the **tyFlow** object by the specified value.

#### **INFO:**

By default, the bounding box of a flow encapsulates all of its particle positions. In order to maximize performance, the bounding box does not encapsulate the shape mesh of each particle, merely the particle's volumeless 3D position in space. Due to this optimization, if all of the particles of a flow are positioned outside of a viewport's frustum, the entire flow may be culled from display even if the shape meshes of particles are large enough to overlap the interior of the frustum. By manually inflating the bounding box of a flow to account for the size of its particle meshes, you can ensure that the flow will not be culled from viewport display even if no particle positions are in view.

#### Miscellaneous

**Ignore .abc notifications**: Ignores rogue change notifications sent by imported Alembic files.

#### **INFO:**

Max's default Alembic importer sends rogue change notifications to its dependent objects when the time slider is moved. Enabling this setting will ignore those notifications, and prevent **tyFlow** simulations which are dependent on Alembic objects from resetting each time the time slider is moved.

**Expand editor rollout**: some Windows display configurations can result in the **tyFlow** editor rollout being cropped improperly. Enabling this setting and increasing the percentage spinner will increase the width of the editor rollout.

%: the percentage (relative to default width) to increase the editor rollout.

**Note:** This is a sticky setting that will persist across max scene files.

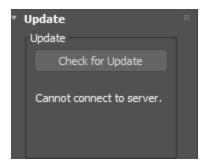
## **About Rollout**

The About rollout contains copyright information, a link to Tyson Ibele's website, as well as third party license information.



# **Update Rollout**

The Update rollout gives you the ability to check online to see if the installed version of **tyFlow** is older than the latest available version.



## **Operators**

Each of these operators will be covered in the following pages: - You can access them from the index system, or on the first page of this helpfile is a image mapped quick link menu where you can click on the operator name to go to that section of the help file.



tyFlow operators are what create and drive particles in a simulation.

## **Operator Activation**

After an operator has been added to a flow, it can be enabled/disabled by clicking on its icon.

## **Operator Settings**

The settings for a particular operator can be accessed in the subrollout on the right hand side of the **tyFlow** editor. An operator must be selected before its settings will be revealed.

## **Common Settings**

Many operators contain common settings. These settings pertain to operator timing controls, as well as test condition controls.

#### Filters Rollout (common Rollout)

The Filters rollout allows you to control which particles will be affected by an operator.



**Enable filters**: controls whether particle filtering will be enabled for the operator.

**AND/OR**: controls the test condition for multiple filters. "And" means that a particle must pass all filters in order to be affected by the operator. "Or" means that a particle must pass any filter in order to be affected by the operator.

**Filter list**: the list of filters that will be used to test particle properties.

**Enabled**: controls whether the selected filter will be enabled.

**Property Type**: the property that the selected filter will test.

**Test TRUE if property is:** the property test condition.

#### **Test Value**

**Absolute/Custom Float**: controls whether the test value will be an absolute value, or a value derived from a custom float data channel of the particle being tested.

Value: the absolute test value.

**Channel**: the custom float data channel from which to derive the test value.

**Variation %**: the per-particle percentage of variation to apply.

#### **Custom Float**

**Channel**: the custom float data channel from which to derive the property value.

#### **Simulation Groups**

**Simulation Groups**: controls which particle simulation groups will be used as the property test.

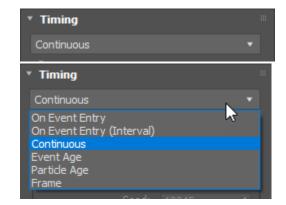
#### Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

#### **Timing Rollout (common Rollout)**

The Timing rollout controls when particles will be affected by an operator.



**Timing type**: controls the timing type of the operator.

**On Event Entry**: particles will only be affected by the operator when they first enter the event.

On Event Entry (Interval): particles will only be affected by the operator if they enter the event within a specified range of frames.

**Continuous**: particles will be continually affected by the operator as long as they stay in the event.

**Event Age**: particles will be continually affected by the operator as long as their lifetime within the event is within a specified range of values.

**Particle Age**: particles will be continually affected by the operator as long as their lifetime within the simulation is within a specified range of values.

**Frames**: particles will be continually affected by the operator as long as the current frame is within a specified range of values.



**Range:** the range of values which will affect various timing types.

Variation: the per-particle amount of variation to apply.

**Seed:** the seed value for all varied parameters.



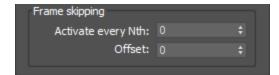
#### Frame exclusions

**Enable frame exclusions**: controls whether the list of frames to exclude will be processed by the operator.

**Frame exclusion list**: the list of frames to exclude from the operator's affect.

#### TIP:

The list may contain multiple frames or frame ranges. Separate frames with commas, and denote ranges of frames with dashes. A valid list may look something like: "0, 10, 20, 35-50, 100" (without quotations).



#### Frame skipping

The frame skipping parameters allow you to control whether an operator will be activated on certain frames, using a modulus operation to easily skip every other frame, or every 3rd frame, etc.

Activate every Nth: controls the divisor of the modulus operation, which takes the form of: if (frame % divisor == 0) {...do work...}. In other words, if the current frame is not evenly divisible by this value, then operator activation on this frame will be disabled. So if you set the value to 2, only every other frame will be activated. If you set it to 3, only every 3rd frame will be activated, etc.

**Offset**: allows you to offset the time value of the modulus operation. An activation value of 2 and an offset value of 0 will activate these frames: 0, 2, 4, 6, 8...etc. An activation value of 2 and an offset value of -1 will activate these frame: 1, 3, 5, 7, 9...etc.

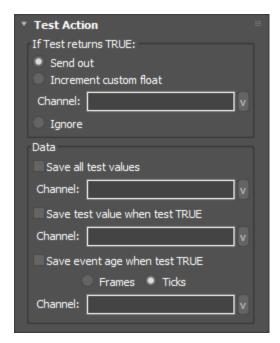
**NOTE:** An activation value of 0 means every whole frame and every substep (that also satisfy the other Timing rollout conditions) will be evaluated. An activation value of 1 means only whole frames will be evaluated (subframes will not be evaluated). An activation value greater than 1 means only every Nth whole frame (where N is the activation value) will be evaluated (subframes will not be evaluated).

#### TIP:

You can speed up the overall evaluation of your flow by using these settings to limit the activation of certain operators. Many operators do not need to evaluate on every timestep or subframe in order to maintain the fidelity of your flow. Using these settings to control frame activation can greatly increase performance in certain situations.

#### **Test Action Rollout (common Rollout)**

The Test Action rollout controls what will happen to particles that satisfy a conditional test within an operator.



**Send Out**: controls whether particles that satisfy the test condition of an operator will be sent to the next event (if the operator has a valid output connection).

**Increment custom float**: controls whether particles that satisfy the test condition of an operator will have the value of one of their custom data channels incremented.

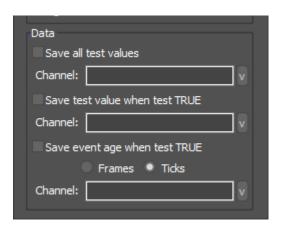
**Channel**: the channel whose value will be incremented.

#### INFO:

Sometimes you may want a particle to satisfy the test conditions of multiple operators before sending it to another event. In that case, you would set the test action of all of the relevant operators to "increment custom float". Then, you would put a Custom Properties operator (whose timing is set to "continuous") at the top of the event's operator stack (and have it reset the value of the appropriate custom data channel to 0 each time step), and a Property Test at the end of the event's operator stack which would be used to check the value of that custom data channel. The Property Test's test action would be set to "send out" and the condition would be set to the desired value of the custom data channel. For example, if you wanted to chain three different operator's test conditions together, you would use the Property Test to test for a custom data channel value of 3 before sending particles to the next event.

With that method in mind you can easily craft more complex test conditions within certain events.

**Ignore**: nothing will happen to particles that satisfy the test condition.



#### **Data**

**Save all test values**: saves any value of the property property to a custom float data channel, regardless of whether or not it satisfied the test condition.

**NOTE:** Enabling the "save all test values" option is a good way to track the properties of particles within tested operators - especially the Property Test operator. You may, for example, save the number of neighbors a particle has from a Property Test, and use that value later as a filter in another operator.

**Save test value when test TRUE**: saves the value of the particle property to a custom float data channel when the value satisfies the test condition.

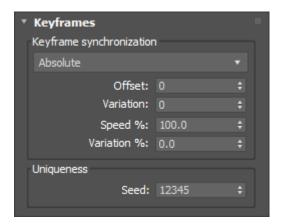
**Save event age when test TRUE**: saves the event age of a particle to a custom float data channel at the time when it satisfies the test condition.

**Frames/Ticks**: controls which unit of time to save the event age value in.

**Channel:** the channel where the value will be saved.

#### **Keyframes Rollout (common Rollout)**

The Keyframes rollout allows to control how animated spinner values will be synchronized - either to the current time or to particle ages.



#### **Keyframe synchronization**

**Absolute**: spinner parameters will be synced to the current time, regardless of particle age.

**Event Age**: spinner parameters will be synced to particle event ages. For example, a particle that has been in an event for 5 frames, will query animated spinner values at frame 5.

**Particle Age:** spinner parameters will be synced to particle ages. For example, a particle that has an age of 5 frames, will query animated spinner values at frame 5.

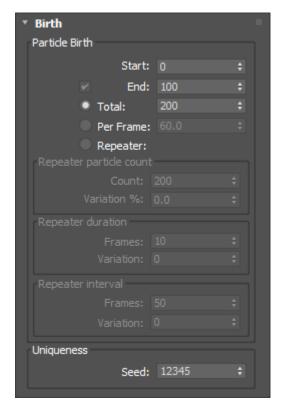
- **Offset**: the per-particle frame offset to apply to the sync method.
- **Variation**: the per-particle variation to apply to the sync method.
- **Speed** %: the rate at which to synchronize keyframes. Lower values will slow animation.
- **Variation** %: the per-particle percentage of variation to apply.

#### Uniqueness

# **Birth operator**



The Birth operator is the simplest operator for creating new particles over time.



**Start/End**: controls the time range in which to birth new particles.

#### NOTE:

When birth mode is not set to "total", you can uncheck "end" in order to have particles be birthed on every frame after the start frame.

**Total:** (Count) controls the total number of particles to birth over the time range.

**Per frame:** (Count) controls the number of particles to birth per frame over the time range.

Repeater:

# Repeater: Repeater particle count

**Count**: controls the number of particles to birth per interval.

**Variation** %: the per-particle percentage of variation to apply.

# Repeater duration

**Frames**: controls the number of active birth frames per interval, within the overall time range.

**Variation**: the per-particle amount of variation to apply.

# Repeater interval

**Frames**: controls the amount of time to wait between birth intervals.

**Variation**: the per-particle amount of variation to apply.

#### TIP

The repeater is useful for repetitive particle birth sequences. For example, imagine you wanted to shoot 50 particles into the scene, every 20 frames. You would enable the repeater and set *count* to 50, *duration* to 1 and *interval* to 20. Using the repeater avoids having to setup multiple birth operators to achieve the same outcome.

# **Birth Burst operator**



The Birth Burst operator allows you to define multiple unique birth events, aka "bursts".



#### Particle Birth Bursts:

**Birth burst list**: the list of all active birth bursts.

# **Burst Settings**

**Start/End**: controls the time range in which the selected burst will birth new particles.

**Total:** (Count) controls the total number of particles the burst will create over the time range.

**Frame (Count)** controls the number of particles the burst will create per frame over the time range.

# Simulation groups

**Simulation groups**: controls which particle simulation groups will be assigned to the particles in the burst.

# **Export groups**

**Export groups**: controls which particle export groups will be assigned to the particles in the burst.

### **Custom float**

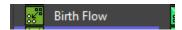
**Value**: the custom float value to assign to the particles in the burst.

**Channel**: the custom float data channel the value will be assigned to.

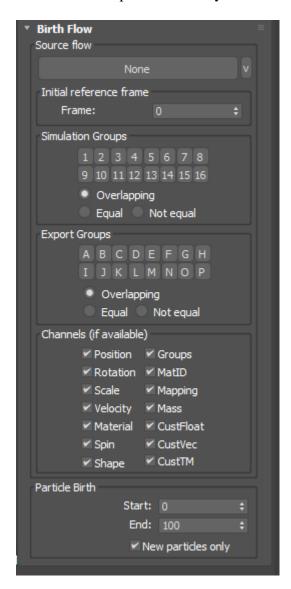
#### **TIP**

The Birth Burst operator is a more efficient alternative to creating multiple unique birth operators in order to birth distinct groups of particles.

# **Birth Flow operator**



The Birth Flow operator allows you to birth new particles that are copies of particles from another flow.



#### Source flow

• **Flow object**: the **tyFlow** object whose particles will be copied.

### Initial reference frame

**Frame**: the initial frame of reference that the input flow object will be evaluated at.

### Simulation Groups

Controls which particle simulation groups will be read from the input flow object.

# Channels (If available)

Controls which particle data channels to copy from the input flow object's particles.

**NOTE:** Birth Flow cannot import bindings/cloth/PhysX data/actor-dependencies/etc from other flows. Only data from the selected channels will carry over.

#### Particle birth

**Start/End**: controls the time range in which to birth new particles, copied from the input flow object.

**New particles only**: controls whether only new particles will be birthed each frame. A particle is considered new if its input ID has not been processed on a previous frame.

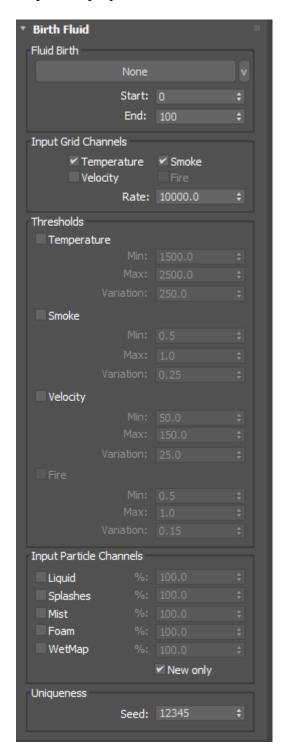
#### **TIP**

Birth flow operators are useful for optimizing the initial states of flows. For example: imagine a flow that computes a voronoi fracture on 1000 particle meshes at frame 0. Every time a property of that flow is changed, the flow's simulation will reset and the voronoi fracture will be recomputed in the process. To avoid having to wait for the fractures to initialize each time the flow property is changed, a secondary flow could be created that uses a birth flow operator in order to copy the fractured particles from the first. Each time the second flow's properties change, the voronoi fracture of the input flow **will not** have to be recomputed as well. And each time the first flow's settings change, the second flow will be automatically updated in the process. In this way, flows can be connected together in order to minimize recomputation costs of complex initial state parameters.

# **Birth Fluid operator**



The Birth Fluid operator can be used to birth particles inside of a (Phoenix FD) fluid grid based on grid voxel and particle properties.



### Fluid Birth

**Fluid object**: the input fluid grid object.

**Start/End**: controls the time range in which to birth new particles.

### Input grid channels

**Temperature**: when enabled, grid cells with a temperature value above 300 will be able to spawn particles.

**Smoke**: when enabled, grid cells with a smoke value above 0 will be able to spawn particles.

**Velocity**: when enabled, grid cells with a velocity value above 0 will be able to spawn particles.

**Fire**: when enabled, grid cells with a fire value above 0 will be able to spawn particles (FumeFX only).

**NOTE:** The higher the temperature/smoke value per cell, the higher the probability that a particle will be spawned in the cell.

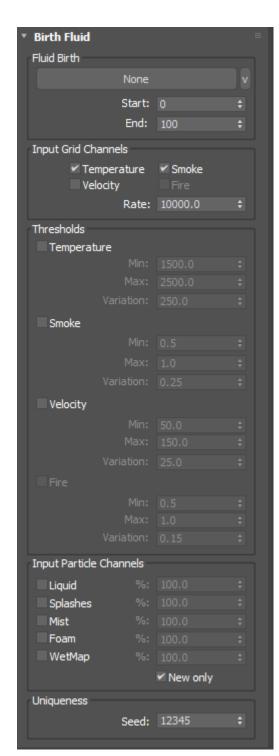
**Rate**: The theoretical rate of particles births that can happen per simulation step, given a per-cell probability of 1. The actual rate of particle births will be lower, and the birth probability per-cell is dependent on the total smoke and temperature values of each cell.

#### **Thresholds**

Enabling these channels will override the default cell spawn probabilities.

**Temperature/Smoke/Velocity/Fire**: the threshold override channels.

**Min/Max/Variation**: the range of values for each selected channel that will control how many particles are spawned in each cell.



# **Input Particle Channels**

PhoenixFD grids can contain both voxels and particles. To birth particles that will be copies of PhoenixFD grid particles, enable the relevant particle channels.

**Channels**: enabling these channels will birth copies of particles contained within them, if such particles are available in the cache of the input object.

**Channel percentages**: controls the percentage of copied input particles to birth.

**New only**: controls whether only new particles will be birthed each frame. A particle is considered new if its input ID has not been processed on a previous frame.

# Uniqueness

# **Birth Intersections operator**



The Birth Intersections operator allows you to birth new particles at places where the geometry of input objects intersects.



#### **Birth Intersections**

**Start/End**: controls the time range in which to birth new particles at intersection points.

**Threshold**: affects the density of particles birthed on two intersecting faces by multiplying the density value by the ratio between this value and the length of the overlap between the intersecting faces.

**Density**: controls the base number of particles to birth over each intersection, prior to adjustments made relative to the threshold value.

**Variation** %: the per-particle percentage of variation to apply.

#### **INFO:**

The number of particles birthed on any given intersecting face can be estimated as:

(density) x (length of intersection overlap / threshold).

#### Mode

A > A: this mode will compute intersections between objects in Geometry List A.

A > B: this mode will compute intersections between objects in Geometry List A and Geometry List B. Intersections between objects in the same group (A > A or B > B) will not be computed.

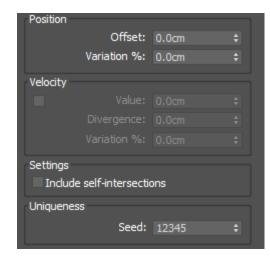
#### **Position**

**Offset**: controls the amount of offset along face normals to position the birthed particles.

**Variation** %: the per-particle percentage of variation to apply.

# Velocity

**Enable velocity**: controls whether velocity along face normals is added to birthed particles.



**Velocity value**: the amount of velocity along face normals to add to birthed particles.

**Divergence**: the angle of divergence applied to velocity vectors.

Variation %: the per-particle percentage of variation to apply.

# Settings

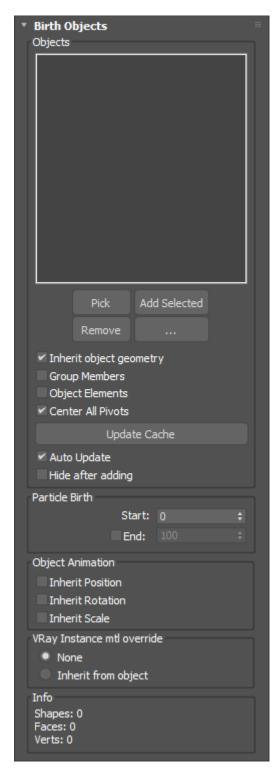
• **Include self-intersections**: controls whether self-intersections will be computed for objects.

# Uniqueness

# **Birth Objects operator**



The Birth Objects operator allows you to convert scene objects into particles.



# **Objects**

**Object list**: the list of input objects which will be converted into particles.

**Inherit object geometry**: controls whether the meshes of input objects will be assigned to corresponding particles.

**Note:** When "inherit object geometry" is disabled, only object transforms will be transferred to birthed particles. This can provide a significant speedup if the input objects have high resolution meshes, but you do not need those meshes to be assigned to the new particles (for example, if you want to align particles to objects but don't need those particles to inherit the meshes).

**Group members**: if an input object is a group head, enabling this setting will convert its constituent members into particles.

**Object elements**: if an input object mesh has multiple subelements, enabling this setting will split them into multiple particles.

**Center all pivots**: centers the pivots of extracted meshes.

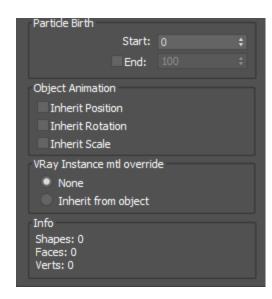
**Update Cache**: click this to manually update the internal cache which holds mesh data for all input objects.

**Auto update**: controls whether changes to input objects will automatically update the operator's internal mesh cache.

**Hide after adding**: controls whether objects will be hidden in the scene after adding them to the listbox.

**Note:** The "hide after adding" setting is a sticky setting (remembered between operators and Max sessions) that will remain checked/unchecked depending on which state you leave it in last.

#### **Birth Objects Operator Continued**



### Particle Birth

**Start**: the start frame at which particles will be birthed.

**End enabled**: controls whether particles should be birthed over a range of frames.

**End**: the end frame at which particles will be birthed.

### **Object Animation**

**Note:** If particles leave the event which contains the Birth Objects operator, their animation will no longer be updated by the Birth Objects operator.

**Inherit Position**: particles will inherit the position changes of the scene object they reference.

**Inherit Rotation**: particles will inherit the rotation changes of the scene object they reference.

**Inherit Scale**: particles will inherit the scale changes of the scene object they reference.

# VRay Instance mtl override

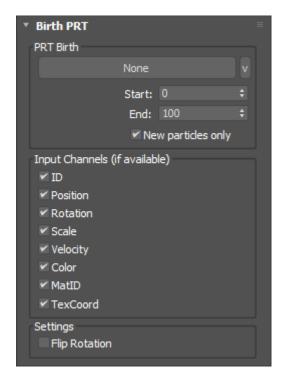
**None**: no material override will be assigned to render instances of the born particles.

**Inherit from object**: the material override for render instances of born particles will be taken from the scene object they are referencing.

# **Birth PRT operator**



The Birth PRT operator allows you to birth new particles that are copies of particles from a PRTLoader object.



### PRT birth

**PRT object**: the input object that contains PRT particles.

**Start/End**: controls the time range in which to birth new particles.

**New particles only:** controls whether only new particles will be birthed each frame. A particle is considered new if its input ID has not been processed on a previous frame, or its age is 0.

# Input channels

**Channels**: controls which PRT data channels to copy into to birthed particles.

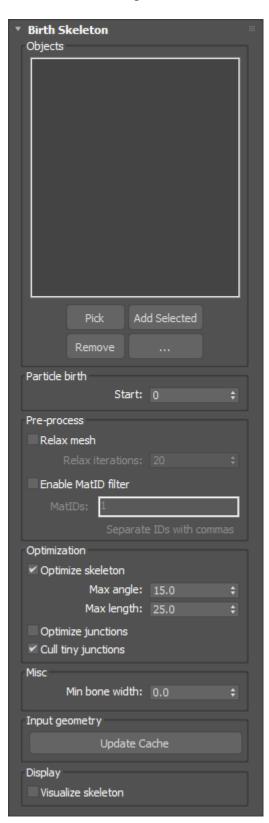
# Settings

**Flip rotation**: flips the w-component of quaternions loaded from PRT data.

# **Birth Skeleton operator**



The Birth Skeleton operator can be used to extract the curve skeleton from input meshes.



#### INFO:

The Birth Skeleton operator extracts the curve skeleton from input meshes, and converts the segments of bones within the extracted skeleton into particles. Using this method, you can easily rig complex meshes like trees and foliage which would otherwise be extremely difficult and time-consuming to manually rig.

#### TIP:

Within the context of this operator, a "segment" is a single line used to compose a bone. A "bone" is a curve composed of one or more "segments". A "skeleton" is the set of all "bones" extracted from a mesh. Keep in mind that the skeleton extracted from a mesh is not a typical anatomical skeleton (ie, this operator is not meant to extract bones for the purpose of character rigging), but a curve skeleton (ie, the algorithm will attempt to convert tube-like sections of a mesh into segmented curves). For this reason, the operator is ideal for extracting bones from tube-like meshes (that's why it works great for plants, foliage, tentacles, etc), but it doesn't work very well for meshes that do not have many long, protruding parts.

# Objects

**Object list**: the list of input objects whose skeleton will be extracted.

#### Particle birth

**Start**: controls the frame at which to birth new particles.

#### **Pre-process**

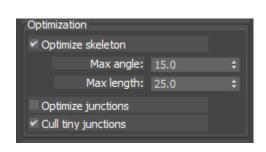
**Relax mesh**: applies laplacian relaxation to the input mesh, in order to smooth out surface details prior to skeletonization.

**Relax iterations**: the number of relax iterations to apply to the input mesh. The higher the iterations, the smoother the mesh will be.

**Enable MatID filter**: when enabled, only faces matching the listed material ID values will be skeletonized.

**MatID**: the list of face material IDs to skeletonize.

#### **Birth Skeleton Operator Continued**



#### TIP:

Input meshes with a lot of small, bumpy surface details can cause artifacts to appear in the resulting skeleton (junctions with unnecessary bones pointing in random directions, bones not properly following the overall surface curvature, etc). Performing pre-process relaxation (or manually relaxing the geometry of the mesh yourself) can help to improve the result of the skeletonization process. An ideal mesh for skeletonization is one that is perfectly smooth and composed only of long, tube-like structures.

#### TIP:

Using the material ID filter is an easy way to exclude parts of a mesh (like leaves, small twigs, etc) from skeletonization.

# Optimization

**Optimize skeleton**: controls whether or not the raw extracted skeleton will be optimized in various ways.

**Note:** The max angle and max length settings control how individual bones will be optimized, by merging/reducing their segments.

**Max angle**: connected segments whose angle is below this threshold will be merged.

**Max length**: segments below the max angle threshold will only be merged if their combined length is below this value.

**Optimize junctions**: in skeletons with clear junctions (places where the endpoint of a bone touches the midpoint segment of another bone), segments will be merged at junction points.

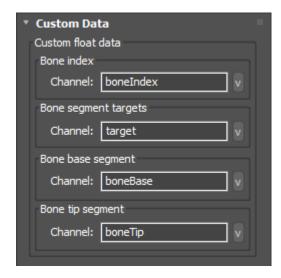
**Note:** Turning "optimize junctions" on may cause endpoints of bones which touch midpoints segments of other bones to lose contact with those midpoint segments, if those midpoint segments are optimized away. Keeping this setting off will maximize junction contact in skeletons extracted from meshes with clean, contiguous topology.

**Cull tiny junctions**: single-segment bones generated at endpoint junctions of multi-segment bones will be culled.

**Note:** The skeletonization algorithm has a tendency to create junctions with multiple single-segment bones pointing outwards at seemingly-random angles, in meshes that aren't composed of perfectly smooth tube-like structures. This setting uses a fairly robust heuristic to minimize the creation of those bones.

#### **Birth Skeleton Operator Continued**







#### Misc

**Min bone width**: specified the minimum width of generated particles.

#### Custom float data

#### **Bone index**

Stores the bone index of each segment into a channel within the corresponding particle.

**Channel**: the custom data float channel where the bone index value will be stored.

#### **Bone segment targets**

Bone segments each have two endpoints (the start and end of each segment line), and bones are composed of multiple segments ordered from base to tip. Thus, the target of each segment is the segment before it (excluding the first segment in a bone, which has no target). Thus, the bone segment target of a corresponding particle will be the particle ID generated before it, as each bone is processed.

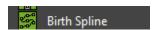
**Channel**: the custom data float channel where the bone segment target value will be stored.

### **Display**

**Visualize skeleton**: draws the raw bone segment lines in the viewport, on the frame they are generated.

**Note:** Each bone is given a random color, and the brightness of the color applied to each bone segment increases from base to tip. Thus, you can see the directionality of each bone in the viewport by looking at the direction of the color gradient for each bone when visualization is enabled. Bone directionality is not controllable, and is determined internally using an algorithm that examines bone radius, proximity, and direction to nearby bones.

# **Birth Spline operator**



The Birth Spline operator can be used to birth new particles on the curves of splines.



### **Splines**

• **Spline list**: the list of input splines whose curves will be used to birth particles.

#### Particle birth

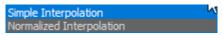


**Percent Along Spline**: particles will be birthed at intervals along splines based on a percentage of their total length.

**Distance Along Spline**: particles will be birthed at intervals along splines based on a distance along their total length.

**Knots**: particles will be birthed on spline knots.

**Interpolated Knots**: particles will be birthed on implicit spline knots, whose interpolated locations are based on the spline's step and optimization settings.



**Simple Interpolation**: percents and distances along the splines will be relative to sub segment lengths and knot spacing.

**Normalized Interpolation**: percents and distances along the splines will be relative to the normalized spline length, ignoring segments lengths and knot spacing.

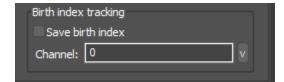
**Start/End**: controls the time range in which to birth new particles.

**Percent** %: the percent value used in "percent along spline" mode.

**Distance**: the distance value used in "distance along spline" mode.

**Relative to scale**: the distance value will be normalized against the magnitude of the spline object's scale vector, so that generated points will be the same distance apart in world space.

#### **Birth Spline Operator Continued**



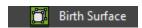
# Birth index tracking

**Save birth index**: controls whether the 0-based index of the spline in the input object list that a particle was birthed on is saved to a the particle's custom data channel.

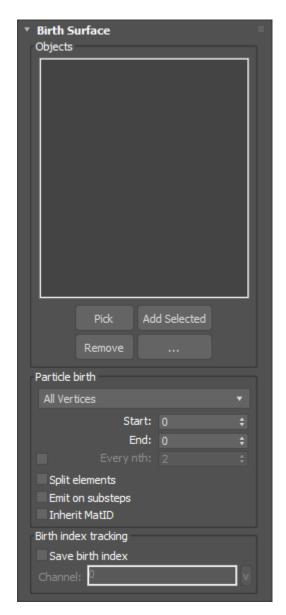
**Note:** The birth index value will be additionally offset by the number of subsplines the input objects have. So, each particle will be given a unique birth index relative to both the input object and its corresponding subspline count. So if two objects are in the list and each object has 1 subsplines, the particle birthed on the first subspline will have an index of 0, and the particle birthed on the last subspline will have an index of 1. If there are two objects in the list, and the first object has 2 subsplines and the second object has 5 subsplines, the particle birthed on the first subspline will have an index of 0, and the particle birthed on the last subspline will have an index of 6.

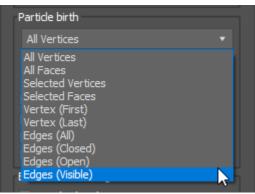
**Channel**: sets the data channel the birth index will be saved to

# **Birth Surface operator**



The Birth Surface operator can be used to birth new particles on specific input mesh features, such as vertices, face centers, etc.





**Note:** Instead of scattering particles randomly on surface features, particles will be sequentially birthed on surface features in the order that the features are indexed. For example, when birthing particles on surface vertices, particle 1 will birth at the location of vertex 1, particle 2 will birth at the location of vertex 2, etc.

# **Objects**

**Object list**: the list of input objects whose surfaces will be used to birth particles.

#### Particle birth

**Surface feature type**: the surface feature on which particles will be birthed.

**Start/End**: controls the time range in which to birth new particles.

**Enable every nth**: when enabled, particles will be birthed at regular frame intervals, instead of at every frame.

**Nth value**: the interval value at which to birth particles.

**Split elements**: controls whether the internal spawn parent value of particles will be incremented depending on which element of a surface the particle is birthed on.

**Emit on substeps**: controls whether particles will be birthed on sub steps of the simulation, or only on whole frames.

**Inherit matID**: controls whether the material ID property of the underlying surface will be copied to the appropriate particle data channel.

#### INFO:

When a particle is birthed on a surface, its internal spawn parent value is set to the index of that surface in the object list, relative to the starting birthID of particles in the event. This means that all particles on the same surface are considered siblings of each other. Their sibling relationship can have an effect on whether they will consider each other binding candidates, among other things.

Sometimes you may not want all particles birthed on surface features to be siblings with each other. For example, if your input object is a shape with multiple spline sub-elements, you would not want particles birthed on separate sub-elements to be siblings of each other. By enabling "split elements", the spawn parent value of birthed particles will not only be incremented for each object in the input list, but also for each sub element of each object. This will create proper sibling relationships between particles birthed on sub-elements of input surfaces, and make it easier to do things like convert input sub-element splines into separate constraint networks.



# Birth index tracking

**Save birth index**: controls whether the 0-based index of the surface in the input object list that a particle was birthed on is saved to a the particle's custom data channel

**Note:** The birth index value will be affected by whether or not "split elements" is enabled.

**Channel**: sets the data channel the birth index will be saved to.

# **Birth Voxels operator**

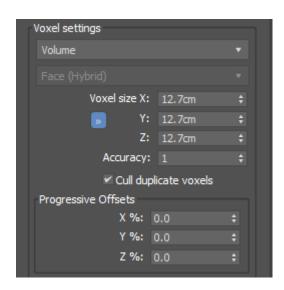


The Birth Voxels operator allows you to birth particles inside the volume or surface of scene objects, at uniform spatial intervals.

#### **Top Part of Rollout**



#### Middle Part of Rollout



# Voxel source objects

**Object list**: the list of input objects whose volumes will be used to birth particles.

#### Particle birth

**Start/End**: controls the time range in which to birth new particles.

**Enable every nth**: when enabled, particles will be birthed at regular frame intervals, instead of at every frame.

**Nth value**: the interval value at which to birth particles.

# **Voxel settings**

**Mode**: controls where particles will be birthed, relative to the surface or volume of a particular input object.

**Sample type**: controls the quality of samples used to determine surface proximity, when birthing particles in "Surface" mode.

**Voxel size X/Y/Z**: the size of individual voxels, in which particles will be birthed.

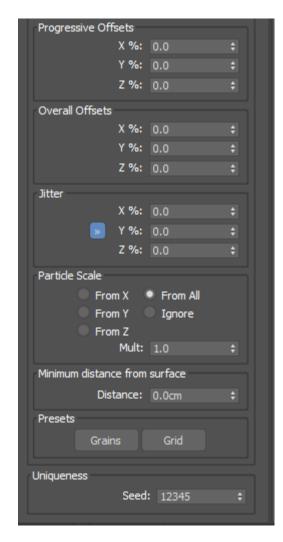
**Accuracy**: controls the accuracy of the raycaster used to compute information about whether a particle is inside or outside of an object's volume. Increase this value for meshes that are self-intersecting or have holes in them.

#### INFO:

If an input object's surface is not closed, or is selfintersecting, increasing the accuracy value can improve results and reduce artifacts.

**Cull duplicate voxels**: when voxelizing multiple input objects, enabling this setting will delete any overlapping particles generated between the objects.

# Birth Voxels Operator Continued Bottom Part of Rollout



#### **Progressive offsets**

**X/Y/Z** %: the percent of progressive position offset applied to each particle, relative to the overall size of the voxels.

#### **Overall offsets**

**X/Y/Z** %: the percent of overall position offset applied to each particle, relative to the overall size of the voxels.

#### **Jitter**

**X/Y/Z** %: the percent of overall position jitter applied to each particle, relative to the overall size of the voxels.

#### **Particle Scale**

**Scale type**: controls which size value particle will derive their scale from.

Mult: a global scale multiplier applied to all particles.

#### Minimum distance from surface

**Distance**: particles within this distance from the surface will be culled.

#### **Presets**

**Grains**: choosing this preset will apply a progressive offset to particles, such that their positions in space will form a tightly-knit overlapping grid.

**Grid**: choosing this preset will remove all offsets from particles, such that their positions in space will form a uniform, aligned grid.

# Uniqueness

# **Array operator**



▶ Filters

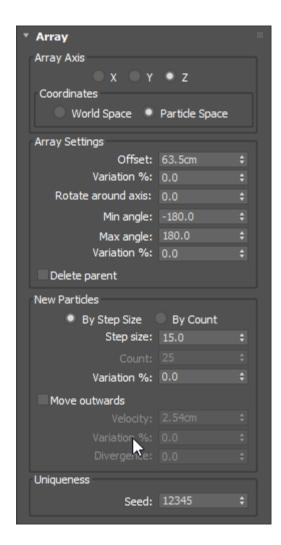
The Array operator allows you to spawn particles in circular patterns.



No help on this rollout at this time



No help on the rollout at this time



# **Array Axis**

**X/Y/Z**: controls the axis around which particles will be spawned.

### Coordinates

**World Space**: the selected axis will be relative to world-space.

Particle Space: the selected axis will be relative to the current particle's transform.

# **Array Settings**

**Offset**: the distance between the current particle and spawned particles.

**Variation** %: the per-particle percentage of variation to apply.

**Rotate around axis:** controls how much rotational offset to apply to all spawned particles, around the array axis.

Min/Max angle: controls the size of the circular arc around the array axis that particles will be spawned within.

**Variation** %: the per-particle percentage of variation to apply.

**Delete parent**: when enabled, the parent particle of the new array particles will be deleted.

#### **New Particles**

By step size: controls whether particles will be spawned at steps (in degrees) around the array arc.

**By count**: controls whether a static number of particles will be spawned, regardless of the size of the array arc.

#### **Array Operator Continued**



**Step size**: the spawn step size, in degrees.

**Count**: the spawn count.

**Variation** %: the per-particle percentage of variation to apply.

**Move outwards**: when enabled, spawned particle velocities will be affected such that they move outward from the current particle.

**Velocity**: the scale of the outward velocity vector to apply to spawned particles.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: controls the degrees of random divergence to apply to outward velocity vectors

# Uniqueness

# **Branch operator**

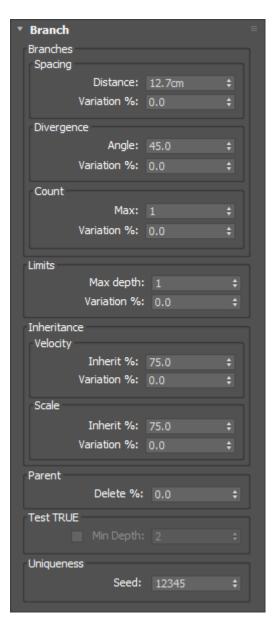


The Branch operator spawns child particles which travel at predictably-divergent angles relative to parent particles. It can be used as the basis for effects like growing frost, lightning, etc.



No help on this rollout at this time

No help on the rollout at this time



#### **Branches**

#### **Spacing**

**Distance**: the minimum distance a parent particle must travel before spawning a child.

**Variation %**: the per-particle percentage of variation to apply.

#### **Divergence**

**Angle**: the angle of divergence applied to spawned particles' velocity trajectory.

**Variation %:** the per-particle percentage of variation to apply.

#### **Count**

**Max**: the maximum number of particles to spawn at each branch event.

**Variation %**: the per-particle percentage of variation to apply.

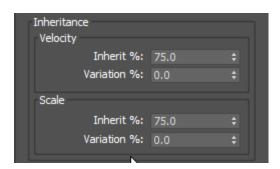
#### Limits

**Max depth**: the maximum number of recursive branches that may be spawned.

Variation %: the per-particle percentage of variation to apply.

**Note:** Increasing max depth increases the number of successive recursive steps the branching algorithm can take. For example, a max depth of 2 means that particles which enter the event may spawn children, and those children may spawn children of their own, but that third group of children (the grand-children of the original particles) may not spawn further children. Increasing the max depth will exponentially increase the total number of particles that will be spawned over time.

#### **Branch Operator Continued**



# Velocity

Inheritance

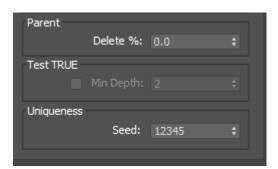
**Inherit** %: the amount of velocity child particles will inherit from parent particles.

**Variation** %: the per-particle percentage of variation to apply.

#### Scale

**Inherit** %: the amount of scale child particles will inherit from parent particles.

**Variation** %: the per-particle percentage of variation to apply



#### **Parent**

No Help at This Time

### **Test TRUE**

**Min depth enable**: controls whether child particles that reach a certain depth will test TRUE for output.

**Min depth value**: sets the minimum depth value required for child particles to satisfy the condition.

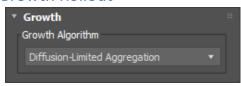
# Uniqueness

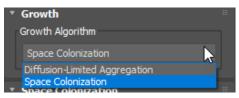
# **Grow operator**



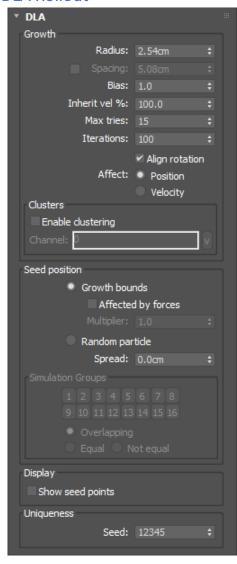
The Grow operator invokes various algorithms to grow intricate patterns of particles.

### **Growth Rollout**





#### **DLA Rollout**



#### **Growth Algorithm**

**Diffusion-Limited Aggregation**: the DLA algorithm grows particles outwards from particles that are surrounded by open space.

**Space Colonization**: the space colonization algorithm grows particles from seed particles that are in close proximity to expansion particles.

#### Growth

**Radius**: controls the size of particles within the DLA algorithm. Increasing this value will increase the overall size of the growth.

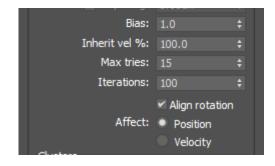
**Note:** As particles are spawned, they are connected to each other's implicit surfaces. The minimum distance between two spawned particles is therefore (**radius x 2**).

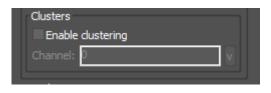
**Enable spacing**: controls whether to use a custom value for the minimum size of empty space required for that region to spawn particles. If this is disabled, the radius value will be used instead.

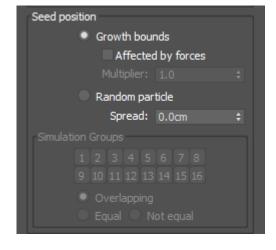
**Spacing value**: the minimum size value of an empty region before it will be able to spawn particles. Increasing this value will result in growths stretching outward, rather than remaining in tighter clumps.

**Note:** Each time the growth algorithm progresses, it chooses a random region adjacent to existing particles and checks that region for overlapping particles. If a particle exists inside the region, spawning inside the region is disabled. The smaller the spacing value, the easier it is for the growth algorithm to find empty regions and therefore progress forward. The larger the region, the more the growth algorithm must stretch out into empty space in order to find an area free of particles.

#### **DLA Rollout Continued**







**Bias**: controls whether the search algorithm will tend to look for valid regions near the center of existing particles, or near the boundary of existing particles. Values less than 1.0 bias towards the center, values greater than 1.0 bias towards the boundary.

**Inherited vel %**: the percentage of velocity child particles will inherit from their parent.

**Max tries**: the maximum number of attempts the algorithm will make to find a valid region of space to spawn a particle, at each growth iteration.

**Iterations**: the number of growth iterations to perform, at each time step of the simulation.

**Align rotation**: aligns the rotation of particles to the vector between themselves and their parent.

**Affect position/velocity**: chooses which particle properties the growth algorithm will affect.

Clusters

**Enable clustering**: controls whether the DLA algorithm will iterate over clusters of particles individually, or over all event particles simultaneously.

**Channel**: the custom particle data channel from which cluster data will be retrieved.

Note: By default, the growth algorithm will process all event particles together. By enabling clustering, you can separate particles into cluster-specific groups, which causes the growth algorithm to process those groups separately. This allows you to have multiple, independent growths forming within the same event.

**Seed position** 

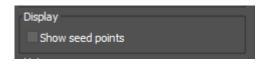
**Growth bounds/random particle**: controls whether starting candidate positions for new growth particles will be located around the bounds of overall growth, or at a random particle within the growth.

**Affected by forces**: controls whether the starting candidate positions for growth particles will be affected by forces.

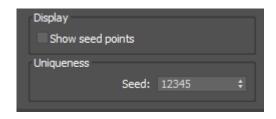
**Multiplier**: the strength multiplier to apply to forces which affect starting candidate positions for growth particles.

**Spread**: the amount of random spread applied to the seed positions.

**Simulation groups**: controls which particle simulation groups the random particle seed position will be derived from.

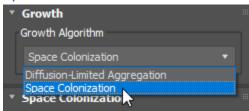


**Show seed points**: displays the resulting seed points in the viewport.



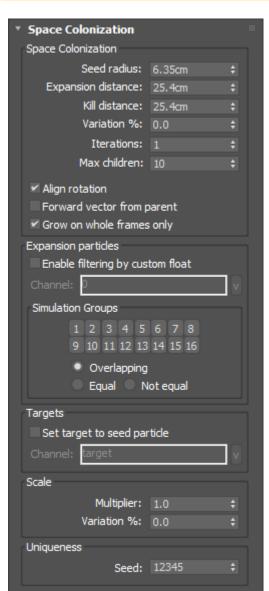
### Uniqueness

# **Space Colonization Rollout**



#### INFO:

Unlike DLA, the space colonization algorithm requires two types of particles in order to run: seed particles and expansion particles. Seed particles are particles which will give birth to new particles during each iteration of the growth cycle. Expansion particles are particles that are in close proximity to seed particles, and that tell seed particles which direction to grow in. A typical setup using this algorithm involves a birth event that generates the initial seed particles, and a birth event which generates the volume of expansion particles the seed particles should grow towards.



**Space Colonization** 

**Seed radius**: controls the size of seed particles.

**Expansion distance**: controls the maximum distance allowed between seed particles and expansion particles. Expansion particles further than this value from any seed particles will not affect growth in the current iteration.

**Kill distance**: controls the maximum distance between expansion particles and seed particles that will cause expansion particles to be deleted after seed particles grow outwards during a growth iteration.

**Variation %:** the per-particle percentage of variation to apply.

**Iterations**: the number of times to repeat the algorithm in a single time step.

**Max children**: the maximum number of children any given seed particle is allowed to grow.

**Align rotation**: aligns the rotation of particles to the vector between themselves and their parent.

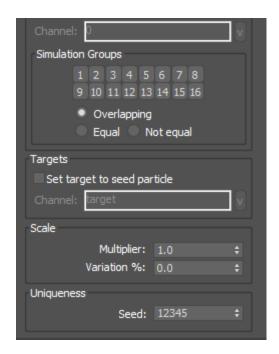
**Grow on whole frames only**: the growth algorithm will only run on whole frames, not sub frames.

### **Expansion particles**

**Enable filtering by custom float**: when enabled, the only particles that will be treated as expansion particles, are those whose custom float data channel value is not equal to 0.

**Channel**: the custom float data channel to use for filtering.

#### **Space Colonization Continued**



**Simulation groups**: controls which particle simulation groups the expansion particles will be derived from.

#### **Targets**

**Set target to seed particle**: each new particle will have its target set to the seed particle it was spawned from.

**Channel**: the custom float data channel to assign the target value to.

#### Scale

**Multiplier**: the multiplier applied to the scale value new particles inherit from their seed parents.

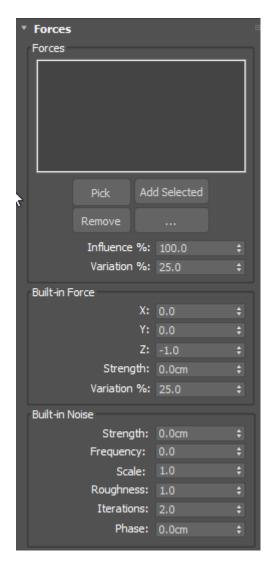
**Variation** %: the per-particle percentage of variation to apply.

#### Uniqueness

### **Forces Rollout**

Forces will affect the direction of growths.

#### **Forces Rollout Top Half**



#### **Forces**

**Force object list**: the list of force objects used to direct the growths.

**Influence** %: the amount of influence forces will have on growths.

**Variation** %: the per-particle percentage of variation to apply.

#### **Built-in Force**

Using the built-in force allows you to apply simple forces to growth particles without having to create force objects in the scene.

**X/Y/Z**: the direction vector of the built-in force.

**Strength**: the strength of the built-in force.

**Variation** %: the per-particle percentage of variation to apply.

#### **Built-in Noise**

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

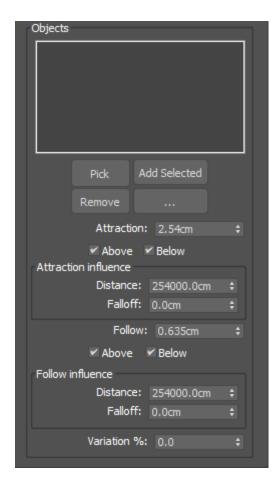
Scale: the scale of the noise.

Roughness: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

**Phase**: the overall evolution offset of the noise pattern.

#### **Forces Rollout Bottom Half**



#### **Objects**

**Object list**: the list of scene objects which will act as growth attractors.

**Attraction**: the amount of attraction forces the scene objects will apply to particles.

**Above/Below**: controls whether attraction forces will be applied to particles above or below the object's surface.

#### **Attraction influence**

The attraction influence extends outwards from nearby objects.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Follow**: the amount of force the scene objects will apply to particles, parallel to the object's surface.

**Above/Below**: controls whether follow forces will be applied to particles above or below the object's surface.

### **Follow influence**

The follow influence extends outwards from nearby objects.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

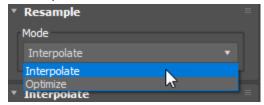
**Variation** %: the per-particle percentage of variation to apply.

# **Resample operator**

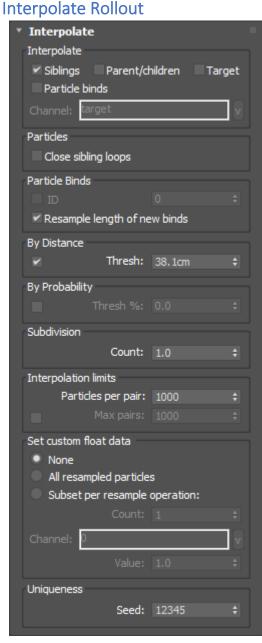


The Resample operator allows you to spawn or delete particles, based on their hierarchical relationship.

# Resample Rollout



# Internalate Pollout



### Mode

**Interpolate**: in this mode, particles will be spawned between matching particle pairs.

**Optimize**: in this mode, particle pairs will be condensed by deleting adjacent particles.

#### **Interpolate**

**Siblings**: particles in an adjacent sibling relationship will be resampled.

**Parent/children**: particles in a parent/child relationship will be resampled.

**Target**: particles and their target (if it exists) will be resampled.

**Particle binds**: controls whether the bindings between particle pairs will be resampled.

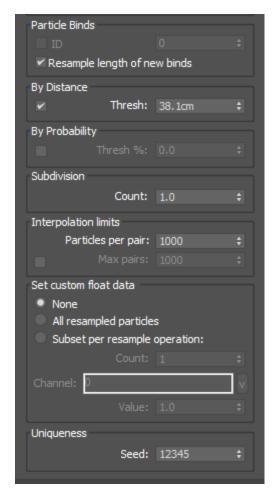
**Target channel**: the custom data float channel from which to get the target particle ID.

**Note:** A particle is a child of another particle if it was spawned from that other particle. Two particles are considered siblings if they were spawned from the same parent particle. Two particles are considered adjacent siblings if they were spawned in order, one immediately before or after the other.

#### **Particles**

**Close sibling loops**: controls whether groups of sibling particles will be turned into closed loops (turn this on to resample particles generated from closed splines).

#### **Interpolate Rollout Continued**



#### **Particle Binds**

**ID**: when enabled, a particle bind's ID must match the ID value in order to be resampled.

**Resample lengths of new binds**: when enabled, new binds (binds created between resampled particles that were previously bound together) will have their reset length set to the interpolated distance between resampled particles.

#### **INFO:**

Example: when "resample lengths of new binds" is enabled, if two particles bound together with a bind length of 5.0 are resampled once (ie, a single new particle is created between them), the resulting two binds created to connect the original particles with the resampled one will each have their rest length set to 2.5. When disabled, they will retain the original bind's rest length of 5.0.

#### By distance

**Enable**: controls whether resampling will occur when two particle pairs are separated by a certain distance.

**Thresh**: the minimum distance required between two particle pairs before resampling will occur.

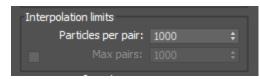
#### By probability

**Enable**: controls whether resampling will occur between particle pairs on a random probability.

**Thresh**: the random probability that two particle pairs will be resampled.

#### **Subdivision**

**Count**: the number of new particles spawned during a resampling of particle pairs.



#### **Interpolation Limits**

**Particles per pair**: limits the total number of new particles that can be spawned between any two matching pairs (sibling or parent/child pairs)

**Enable max pairs**: limits the total number of new particles that can be spawned, per time step.

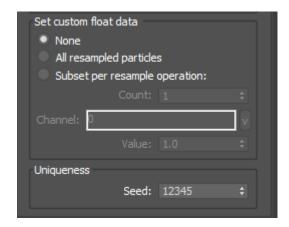
**Max pairs**: the maximum number of particle pairs that can be resampled, per time step.

#### TIP:

When "particle binds" is enabled, if pair particles A and B are resampled such that particle C is spawned between them, any bindings between A and B will be resampled such that A will be bound to C, and C will be bound to B.

$$(A-B) > (A-C-B)$$

#### **Interpolate Rollout Continued**



#### Set custom float data

**None**: no resampled particles will have a custom float value assigned to them.

**All**: all resampled particles will have a custom float value assigned to them.

**Subset per resample operation**: a certain number of resampled particles, per resample operation, will have a custom float value assigned to them.

**Channel**: the custom float data channel to assign a value to.

Value: the value to assign to the custom float data channel.

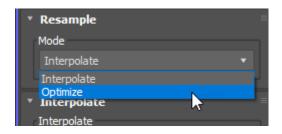
#### TIP:

Use the custom float data settings to mark certain resampled particles, for further processing later.

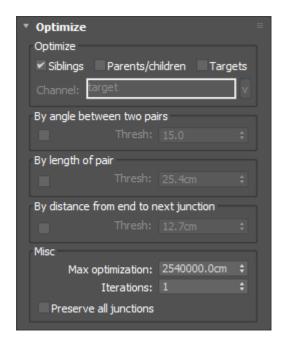
#### Uniqueness

# **Optimize Rollout**

#### To Access Optimize Rollout



#### **Optimize Rollout:**



#### Note:

Certain optimizations can result in more particles becoming optimization candidates after the algorithm completes one iteration. Increasing the iteration count will optimize those new candidates in the same time step.

#### **Optimize**

**Siblings**: particles in an adjacent sibling relationship will be optimized.

**Parent/children**: particles in a parent/child relationship will be optimized.

**Target**: particles and their target (if it exists) will be optimized.

**Target channel**: the custom data float channel from which to get the target particle ID.

#### By angle between two pairs

**Enable**: controls whether particles will be deleted if the angle between their adjacent pairs is below a certain threshold.

**Thresh**: the angle threshold.

### By length of pair

**Enable**: controls whether particles will be deleted if the distance between them and their adjacent particle is below a certain threshold.

**Thresh**: the length threshold.

#### By distance from end to next junction

**Enable**: controls whether particles adjacent to a single junction particle will be deleted if the distance between them and the junction is below a certain threshold.

**Thresh**: the distance threshold.

#### Misc

**Max optimization**: particles will only be deleted if the length of the resulting pair (the pair formed by both of the deleted particle's adjacent particles) is below this threshold.

**Iterations**: the number of times to re-perform the overall optimize algorithm.

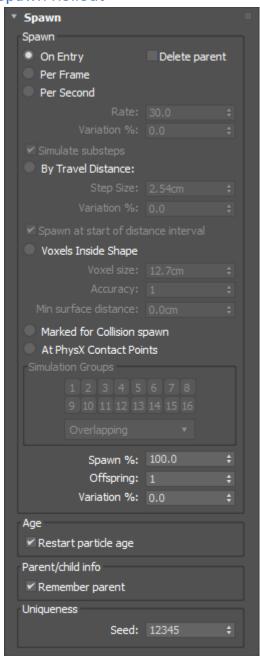
**Preserve all junctions**: junctions (particles with more than two adjacent particles) will not be deleted.

# Spawn operator



The Spawn operator spawns new particles from existing particles.

## Spawn Rollout



### Spawn



Spawns child particles the first time a particle enters the operator

Spawns child particles at a rate per frame over time Spawns child particles at a rate per second over time.



**Delete parent**: controls whether the parent is deleted after children are spawned.

**Rate**: the spawn rate per second.

**Note:** The rate per frame can be calculated as (rate / framerate).

**Variation** %: the per-particle percentage of variation to apply.

**Simulate substeps**: controls whether particle positions will be interpolated along their velocity by a random value, to simulate substep spawning.

## **By Travel Distance:**

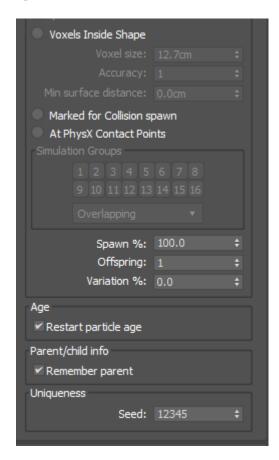
Spawns child particles along the trajectories of their parents, at defined spatial increments.

**Step Size**: the distance a parent particle must travel before spawning a child particle.

**Note:** If a parent particle travels a distance greater than the step size value over a particular step of the simulation, the distance traveled will be properly subdivided such that child particles will be spawned at equal distance intervals.

Variation %: the per-particle percentage of variation to apply

## **Spawn Rollout Continued**



### **Voxels Inside Shape**

Spawns child particles inside voxels within the overall volume of a parent particle's shape mesh.

**Voxel size**: the size of individual voxels, in which particles will be spawned.

**Accuracy**: controls the accuracy of the raycaster used to compute information about whether a particle is inside or outside of a shape mesh's volume.

### INFO:

If a parent particle's shape mesh is not closed, or is selfintersecting, increasing the accuracy value can improve results and reduce artifacts.

**Min surface distance**: particles within this distance from the parent particle's shape mesh will be culled.

## **Marked for Collision Spawn**

Spawns child particles if their parent was marked by a Collision operator as having colliding with a surface at the end of the previous time step.

## **At PhysX Contact Points**

Spawns child particles at contact reports reported by the PhysX solver located on a parent particle's PhysX hull.

**Simulation groups**: controls which particle simulation groups value the PhysX contacts must satisfy, in order to spawn particles.

### INFO:

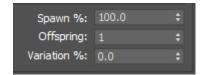
By filtering desired simulation groups, you can control which types of PhysX contacts will spawn particles.

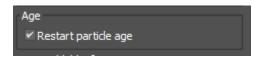
**Spawn %**: the percentage of parent particles that will spawn child particles.

**Offspring**: the number of child particles a particular parent particle will spawn, per time step.

**Variation** %: the per-particle percentage of variation to apply. **Age** 

**Restart particle age**: controls whether spawned particles will have their age reset, or instead inherit the age of their parent.







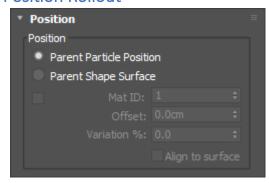


### Parent/child info

**Remember parent**: controls whether the particle will internally remember which particle is its parent. **Uniqueness** 

**Seed**: the seed value for all varied parameters.

## **Position Rollout**



### **Position**

**Parent particle position**: spawned particles will be located at the position of their parent.

**Parent shape surface**: spawned particles will be located at a random position on their parent particle's shape surface.

**Mat ID enabled**: controls whether particles will only be spawned on faces with a specific material ID.

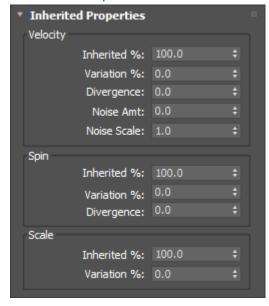
**Mat ID**: the specific material ID of faces particles will be spawned on.

**Offset**: the distance from the parent particle's shape surface, along its face normals, to offset spawned particles.

**Variation** %: the per-particle percentage of variation to apply.

**Align to Surface**: controls whether particles will be aligned to the surface of their parent.

## **Inherited Properties Rollout**



## **Velocity**

**Inherited** %: controls how much velocity child particles will inherit from their parent.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: controls how much child particle velocity vectors will diverge from that of their parent.

**Noise amount**: controls the amount of noise variation that will be applied to child particle velocity vectors.

**Noise scale**: controls the scale of the noise variation applied to child particle velocity vectors.

## **Spin**

**Inherited %**: controls how much spin child particles will inherit from their parent.

**Variation** %: the per-particle percentage of variation to apply.

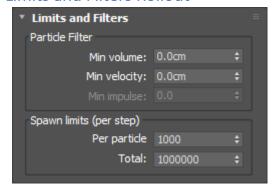
**Divergence**: controls how much child particle spin vectors will diverge from that of their parent.

### Scale

**Inherited %**: controls how much scale child particles will inherit from their parent.

**Variation** %: the per-particle percentage of variation to apply.

## Limits and Filters Rollout



### **Particle Filter**

**Min volume**: parent particles with a shape mesh volume below this value will not spawn children.

**Min velocity**: parent particles with a velocity magnitude below this value will not spawn children.

**Min impulse**: parent particles with PhysX contacts whose impulse is below this threshold will not spawn children.

### **Spawn Limits (per step)**

**Per particle**: limits the total number of children that an individual particle can spawn, per time step.

**Total**: limits the total number of children that can be spawned, per time step.

# **Delete operator**



The Delete operator allows you to delete particles from the simulation.

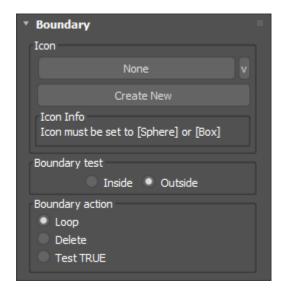
## **Note:**

The Delete operator has no unique parameters. Deletion is instead controlled by parameters in the operator's Timing rollout.

# **Boundary operator**



The Boundary operator allows you to constrain particles to a **tyIcon** volume. Particles outside the volume can be set to loop within the boundary of the icon, be deleted, or test TRUE to satisfy the output condition.



## **Icon**

**Icon object**: the icon object to use as the boundary.

## **Boundary Test**

**Inside**: particles will be tested to see if they are inside the boundary.

**Outside**: particles will be tested to see if they are outside the boundary.

## **Boundary Action**

**Loop**: particles that pass the boundary test will loop to the opposite side of the boundary.

**Delete**: particles that pass the boundary test will be deleted.

**Test TRUE::** particles that pass the boundary test will satisfy the output condition.

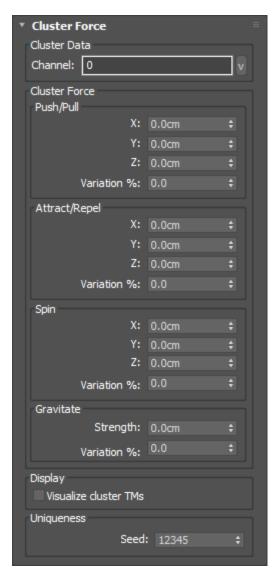
# **Cluster Force operator**



The Cluster Force operator allows you to apply localized forces to clusters of particles. The forces are derived from the eigenvectors of the clusters (equivalent to the transform of the object-oriented bounding box of the clusters).

### INFO:

When using the Cluster Force operator, you should generate the clusters using a Cluster operator set to "LDNP" mode.



### **Cluster Force**

### **Push/Pull**

Forces will push/pull particles along the cluster eigenvectors.

**X/Y/Z**: the strength of the forces along each local eigenvector axis.

**Variation** %: the per-particle percentage of variation to apply.

## Attract/Repel

Forces will attract/repel particles towards/away the cluster eigenvectors.

**X/Y/Z**: the strength of the forces towards/away each local eigenvector axis.

**Variation %:** the per-particle percentage of variation to apply.

## **Spin**

Forces will spin particles around the cluster eigenvectors.

**X/Y/Z**: the strength of the forces around each local eigenvector axis.

**Variation** %: the per-particle percentage of variation to apply.

## Gravitate

Forces will move particles towards/away the cluster centers.

**Strength**: the strength of the local gravity forces.

**Variation** %: the per-particle percentage of variation to apply.

## **Display**

**Visualize cluster TMs**: displays each cluster's eigenvector axis in the viewport

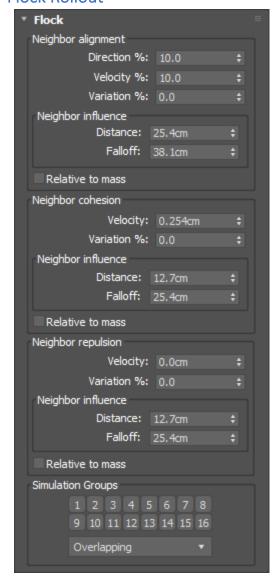
## Uniqueness

# Flock operator



The Flock operator allows you to add boid-style flocking forces to particles.

## Flock Rollout



## **Neighbor Alignment**

**Direction %:** the percentage of cumulative neighbor direction each particle will adopt.

**Velocity** %: the percentage of cumulative neighbor velocity each particle will adopt.

**Variation** %: the per-particle percentage of variation to apply.

## **Neighbor influence**

The neighbor influence extends outwards from the particle's neighbors

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Relative to mass**: the force applied to a particle will be relative to its mass.

## **Neighbor Cohesion**

**Velocity**: the velocity a particle will travel towards its neighbors.

**Variation** %: the per-particle percentage of variation to apply.

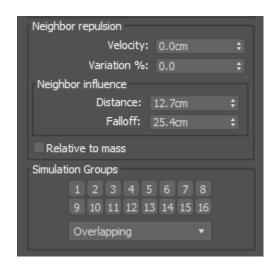
## **Neighbor influence**

The neighbor influence extends outwards from the particle's neighbors

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

### **Flock Rollout Continued**



## **Neighbor Repulsion**

**Velocity**: the velocity a particle will travel away from its neighbors.

Variation %: the per-particle percentage of variation to apply.

## **Neighbor influence**

The neighbor influence extends outwards from the particle's neighbors

**Distance**: particles within this distance will be fully affected.

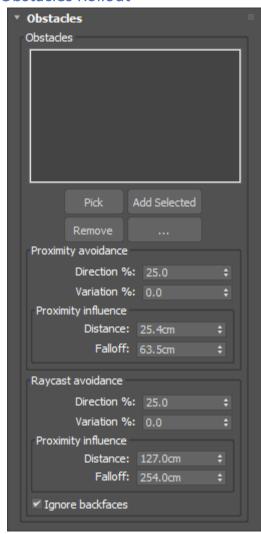
**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Relative to mass**: the force applied to a particle will be relative to its mass.

## **Simulation groups**

• **Simulation groups**: only neighbors with matching simulation groups will be considered.

## **Obstacles Rollout**



### **Obstacles**

Object list: the obstacle objects.

### **Proximity Avoidance**

**Direction %**: the percentage of avoidance direction each particle will adopt.

**Variation %**: the per-particle percentage of variation to apply.

## **Proximity influence**

The proximity influence extends outwards from the surface of nearby obstacles.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

## **Raycast Avoidance**

**Direction %**: the percentage of avoidance direction each particle will adopt.

**Variation** %: the per-particle percentage of variation to apply.

## **Proximity influence**

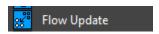
The proximity influence extends outwards from the surface of obstacles hit by raycasts.

**Distance**: particles within this distance will be fully affected.

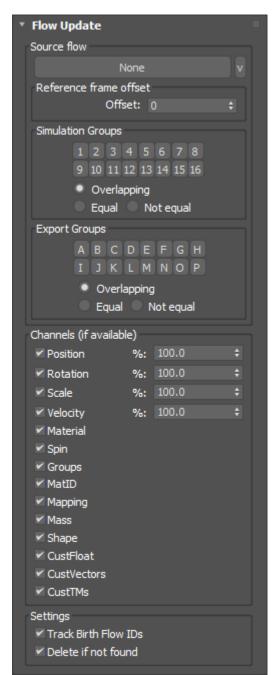
**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Ignore backfaces**: rays that hit backfaces of obstacles will be ignored.

# Flow Update operator



The Flow Update operator allows you to update particle properties based on matching particles within another flow.



### Source flow

**Flow object**: the **tyFlow** object whose particles will be referenced.

**Reference frame offset**: the frame offset that the input flow object will be evaluated at.

**Simulation groups**: controls which particle simulation groups will be read from the input flow object.

**Channels**: controls which particle data channels to copy from the input flow object's particles.

### **Settings**

**Track Birth Flow IDs**: when enabled, particles will be matched using birth IDs previously imported with a Birth Flow operator. When disabled, each particle's own birth ID will be used to find matching particles in the reference flow.

## **INFO:**

If you are using Flow Update in combination with Birth Flow, you should keep "track Birth Flow IDs" enabled. If you are using Flow Update on its own in order to transfer properties from any arbitrary external flow or particle system, you should disable "track Birth Flow IDs".

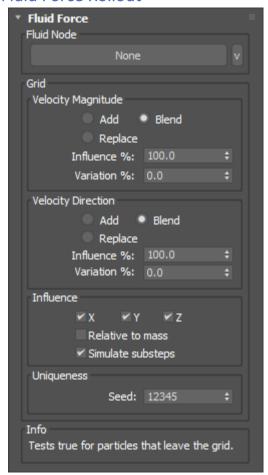
**Delete if not found**: if a particle does not have a matching particle in the input flow object, it will be deleted.

# Fluid Force operator



The Fluid Force operator allows you to apply forces from (Phoenix FD and FumeFX) fluid grids to particles.

### Fluid Force Rollout



#### Fluid Node

**Fluid input object**: the grid object that forces will be loaded from.

### Grid

### **INFO:**

Each particle velocity has a direction and a magnitude. Normally these values are processed as a single vector, but the Fluid Force operator gives you individual controls for both, which allow you to exert fine control over how fluid grids will affect particles.

## **Velocity Magnitude**

These controls affect the overall magnitude of fluid forces applied to particle velocities.

**Add/Blend/Replace**: controls how forces computed by the operator will affect existing particle velocities.

**Influence** %: controls how much influence fluid forces will have on particle velocities.

**Variation** %: the per-particle percentage of variation to apply.

## **Velocity Direction**

These controls affect the overall direction of fluid forces applied to particle velocities.

**Add/Blend/Replace**: controls how forces computed by the operator will affect existing particle velocities.

**Influence** %: controls how much influence fluid forces will have on particle velocities.

**Variation** %: the per-particle percentage of variation to apply.

### **Influence**

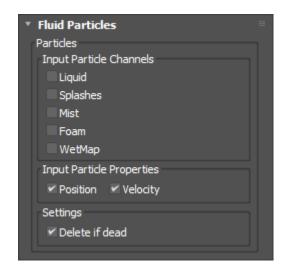
**X/Y/Z**: controls which axis the forces will affect.

**Relative to mass**: controls whether particle forces will be relative to particle masses.

**Simulate substeps**: forces will be interpolated in a way that simulates the addition of forces at smaller simulation substeps.

### **Uniqueness**

## Fluid Particles Rollout



### **Particles**

PhoenixFD liquid grids may also contain particle data. These controls allow you to drive the corresponding particles birthed from a Birth Fluid operator, using the same grid.

## **Input Particle Channels**

**Channels**: the various PhoenixFD particle channels which may be available in the grid.

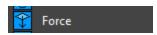
## **Input Particle Properties**

**Position/Velocity**: controls which properties to copy from PhoenixFD particles onto the corresponding particles within the event.

## **Settings**

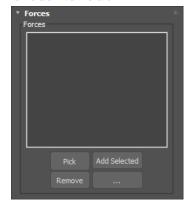
**Delete if dead**: if a particle from the event has no corresponding particle in the fluid grid for the selected input channels, it will be deleted.

# Force operator



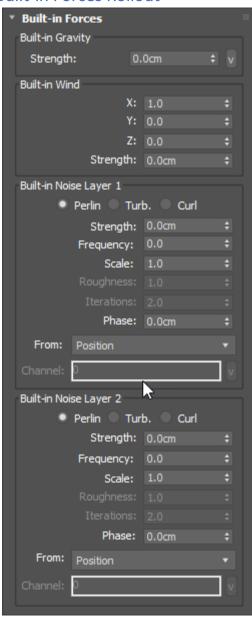
The Force operator allows you to apply various forces to particles.

## **Forces Rollout**



**Force objects**: The input force/spacewarp helpers.

## **Built-in Forces Rollout**



## **Built-in Gravity**

**Strength**: the strength of the built-in gravity to apply to particles.

**Built-in Wind** 

**X/Y/Z**: controls the direction vector of the wind.

**Strength**: controls the strength of the wind.

### **Built-in Noise Layer 1 & 2**

## TIP:

Using both built-in noise layers with different settings can improve the visual complexity of the resulting noise values.

**Noise type**: controls which noise algorithm will be used to calculate forces.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will increase force evolution over time.

**Scale**: the scale of the noise.

**Roughness**: the roughness of the turbulent noise.

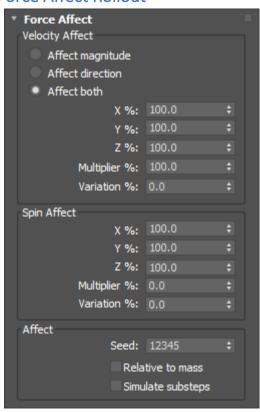
**Iterations**: the number of iterations to use to calculate the turbulent noise.

**Phase**: the overall evolution offset of the noise pattern.

**Source vector**: controls the input vector used to calculate the noise values.

**Channel**: the custom data channel to read custom vectors from.

## **Force Affect Rollout**



### **Velocity Affect**

**Affect magnitude/direction/both**: controls which aspects of the resulting particle velocity the forces will affect.

**X/Y/Z** %: controls the amount of influence forces will have over each axis of the resulting velocity.

**Multiplier** %: an overall multiplier applied to the resulting force values.

**Variation** %: the per-particle percentage of variation to apply.

## **Spin Affect**

**X/Y/Z** %: controls the amount of influence forces will have over each axis of the resulting spin.

**Multiplier** %: an overall multiplier applied to the resulting force values.

**Variation** %: the per-particle percentage of variation to apply.

**Note:** When applying forces to spin values, the influence percentage may need to be very high to get the desired affect, due to the fact that spin is calculated in degrees per second and velocity is calculated in scene units per time step.

### **Affect**

**Seed**: the seed value for all varied parameters.

**Relative to mass**: the force applied to a particle will be relative to its mass.

**Simulate substeps**: forces will be interpolated in a way that simulates the addition of forces at smaller simulation substeps.

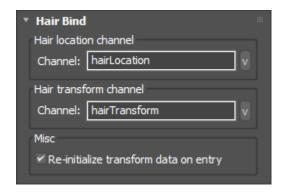
# **Hair Bind operator**



The Hair Bind operator allows you to bind particles to Ornatrix hair strands.

## **INFO:**

The Hair Bind operator requires hair strand location data derived from a Position Hair operator. Make sure you setup a Position Hair operator above the Hair Bind operator in your flow.



### Hair location channel

**Channel**: the custom TM channel from which strand location data (setup in a prior Position Hair operator) will be derived.

## Hair transform channel

**Channel**: the custom TM channel to which local strand transform data will be saved/loaded.

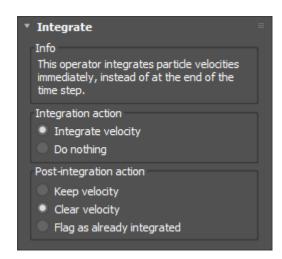
### Misc

**Re-initialize transform data on entry**: when particles enter the event, their transform channel data (their strand-space local transform) will be reset.

# **Integrate operator**



The Integrate operator allows you to immediately integrate particle velocities into particle positions.



### **Integration action**

**Integrate velocity**: adds the velocity of the particle to the position of the particle.

**Do nothing**: no integration action will be taken.

### **Post-integration action**

**Keep velocity**: after the velocity is integrated, the velocity is left unchanged.

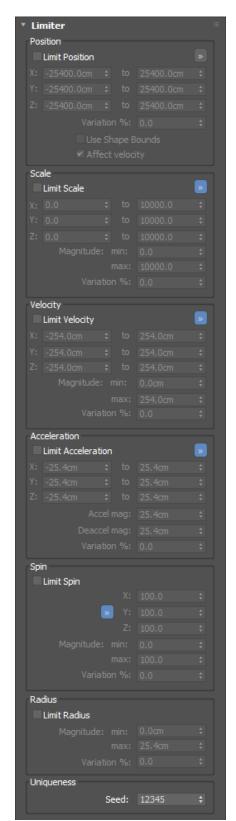
**Clear velocity**: after the velocity is integrated, the velocity is cleared (set to zero).

**Flag as already integrated**: after the velocity is integrated, the velocity is left unchanged but also flagged internally as already integrated, so that it won't be re-integrated at the end of the time step.

# Limiter operator



The Limiter operator can be used to constrain various particle properties.



### **Position**

**Limit position**: controls whether the position values of particles will be clamped.

**X/Y/Z**: sets the min/max allowable values.

**Variation** %: the per-particle percentage of variation to apply.

**Use shape bounds**: considers the bounding box of the particle shape when clamping its position.

**Affect velocity**: controls whether velocity values are cleared for particles whose position is limited.

**Note:** By default, when a particle position is outside the bounds of the limiter values it velocity along the offending axis will be set to 0. This is so that the velocity integrator doesn't move the particle past the limited value during the velocity integration function at the end of the time step. Sometimes this is undesirable behavior, so the ability to disable this has been added.

#### Scale

**Limit scale**: controls whether the scale values of particles will be clamped.

**X/Y/Z**: sets the min/max allowable values.

**Magnitude min/max**: sets the minimum/maximum overall magnitude of particle scale vectors.

**Variation** %: the per-particle percentage of variation to apply.

## Velocity

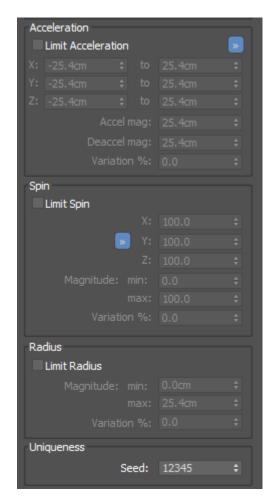
**Limit velocity**: controls whether the velocity values of particles will be clamped.

**X/Y/Z**: sets the min/max allowable values.

**Magnitude min/max**: sets the minimum/maximum overall magnitude of particle velocity vectors.

**Variation** %: the per-particle percentage of variation to apply.

### **Limiter Operator Continued**



### Acceleration

**Limit acceleration**: controls whether the acceleration values of particles will be clamped.

**X/Y/Z**: sets the min/max allowable values.

**Accel mag**: sets the maximum overall magnitude of particle acceleration vectors.

**Deaccel mag**: sets the maximum overall magnitude of particle deacceleration vectors.

**Variation** %: the per-particle percentage of variation to apply.

## Spin

Limit spin: controls whether the spin values of particles will be clamped.

X/Y/Z: sets the min/max allowable values.

**Magnitude min/max**: sets the minimum/maximum overall magnitude of particle spin vectors.

**Variation** %: the per-particle percentage of variation to apply.

### Radius

**Limit radius**: controls whether the radius values of particles will be clamped.

**Magnitude min/max**: sets the minimum/maximum overall magnitude of particle radius values.

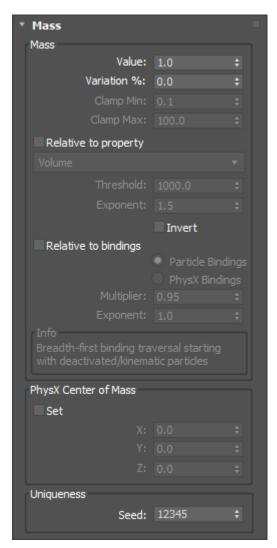
**Variation** %: the per-particle percentage of variation to apply.

## Uniqueness

# **Mass operator**



The Mass operator allows you to control the mass of particles. The mass value of a particle will affect the way it behaves within the bind solver, as well as the PhysX solver.



### **PhysX Center of Mass**

**Set**: controls whether to override the local center of mass offset value on PhysX particles.

**X/Y/Z**: the axis values for the center of mass offset, relative to the local particle transform.

### Mass

**Value**: the initial absolute mass value assigned to particles.

Variation %: the per-particle percentage of variation to apply

**Clamp min**: the minimum allowable mass value.

Clamp max: the maximum allowable mass value.

**Relative to property**: controls whether assigned mass values will be relative to a particle property.

**Threshold:** the target threshold that particle property will be compared to, when assigning mass values.

**Target volume**: the target volume that particle volumes will be compared to, when assigning mass values.

**Exponent**: the exponent that particle volume ratios will be raised to. Larger exponent values will increase the mass disparity between large and small volumes.

**Invert**: controls whether the resulting volume ratio will be inverted.

#### INFO:

When "relative to volume" is on, the following equation is used to calculate particle masses:

mass = pow(volume/(target volume), exponent)

**Relative to bindings**: controls whether masses will be affected by their distance to deactivated/kinematic particles in their binding network.

**Particle bindings**: regular particle bindings will be considered in the breadth-first traversal of the binding network.

**PhysX bindings**: PhysX bindings will be considered in the breadth-first traversal of the binding network.

**Multiplier**: the multiplier applied to masses, relative to their depth in the breadth-first search.

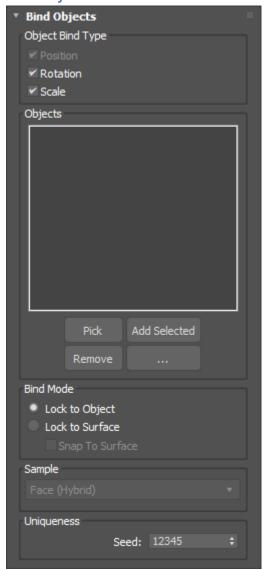
**Exponent**: the exponent applied to the depth-relative multiplier. The larger the exponent, the greater affect the multiplier will have on particles relative to their depth in the breadth-first search

# **Object Bind operator**



The Object Bind operator can be used to bind particles to objects and their surfaces. Various settings allow you to control the strength of the binds across multiple particle data channels (position/rotation/scale), as well as various other surface constraints.

## **Bind Objects Rollout**



### **Object Bind Type**

**Position**: binds particle positions to objects.

**Rotation**: binds particle rotations to objects.

**Scale**: binds particle scales to objects.

## **Objects**

**Object list**: the list of objects to which particles will be bound.

### **Bind Mode**

**Lock to object**: particles will be bound to object transforms.

**Lock to surface**: particles will be bound to object surfaces.

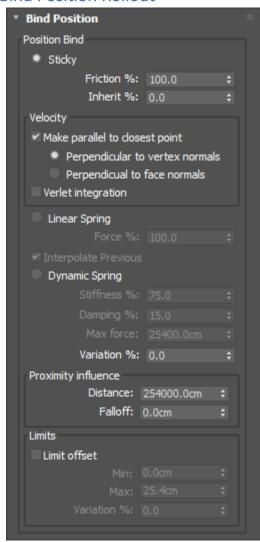
**Snap to surface**: particles bound to surfaces will immediately snap to valid surface locations.

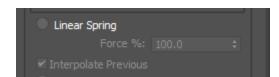
## Sample

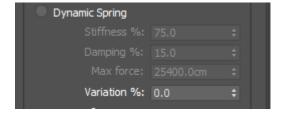
**Sample type**: controls which sampler will be used to determine closest-surface proximities for particles.

## Uniqueness

## **Bind Position Rollout**







### **Position Bind**

### **Sticky**

Sticky bindings constrain particles to consistent distances from their target surface.

**Friction** %: controls the amount of resistance a particle will exert on forces that try to move it along its target surface.

**Inherit** %: controls the amount of velocity particles will inherit from the motion of their target surface.

## **Velocity**

**Make parallel to closest point**: particle velocities will be projected along the nearest point on the object to which they are bound.

**Note:** When bind mode is set to "lock to object", the closest point will be the object's pivot. When bind mode is set to "lock to surface", the closest point will be the closest point on the object's surface.

**Relative to face/vertex normals**: controls which mesh element the projected velocities will be made perpendicular to.

**Verlet integration**: after all bind calculations are completed, the particle velocity will be set to the vector between the final particle position and the start particle position.

**Linear Spring** 

Linear Spring bindings modify particle velocities such that they point directly at their target surface locations.

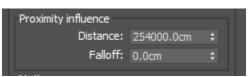
**Force** %: the force of the resulting velocity.

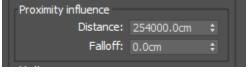
**Interpolate previous**: controls whether particle velocities will be interpolated with their previous values, based on the force multiplier.

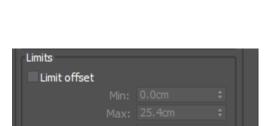
## **Dynamic Spring**

Dynamic Spring bindings modify particle velocities such that target surface vectors are added to their current velocity vectors.

**Stiffness %:** the strength of the target surface vector.







**Damping %:** the amount of damping to apply to existing velocities, prior to the addition of the target surface vector.

**Max force**: the maximum magnitude of target surface vectors.

**Variation** %: the per-particle percentage of variation to apply. **Proximity influence** 

The proximity influence extends outwards from the bind object's closest point.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

### Limits

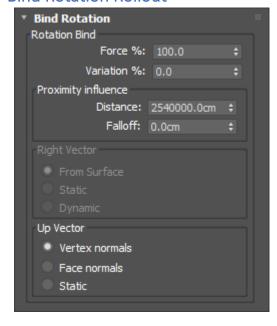
**Limit offset**: limits the distance between a particle's bind target and the actual distance to the nearest point on the surface.

Min: the minimum allowable distance between a bind target and the nearest point on the surface.

Max: the maximum allowable distance between a bind target and the nearest point on the surface.

**Variation %:** the per-particle percentage of variation to apply.

## **Bind Rotation Rollout**



Force %: the amount of force used to constrain particle rotations to their initial orientation offset from the surface.

**Variation** %: the per-particle percentage of variation to apply.

## **Proximity influence**

The proximity influence extends outwards from the bind object's closest point.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

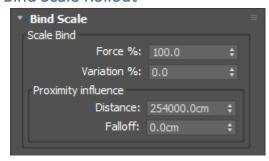
## **Right Vector**

**Right vector type**: the location relative to the surface or the world from which right vectors of the binding orientation matrix will be derived.

## Up Vector

**Up vector type**: the location relative to the surface or the world from which up vectors of the binding orientation matrix will be derived.

## **Bind Scale Rollout**



## Bind Scale

**Force** %: the amount of force used to constrain particle scales to their initial scale offset from the surface.

**Variation** %: the per-particle percentage of variation to apply.

## **Proximity influence**

The proximity influence extends outwards from the bind object's closest point.

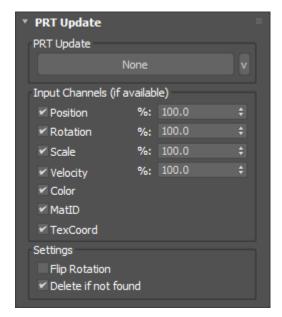
**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

# **PRT** Update operator



The PRT Update operator allows you to update particle properties based on matching particles within a PRTLoader object.



## **PRT Update**

**PRT object**: the PRTLoader object whose particles will be referenced.

## **Input Channels**

**Channels**: controls which particle data channels to copy from the input PRTLoader object's particles.

## **Settings**

**Flip rotation**: flips the w-component of quaternions loaded from PRT data.

**Delete if not found**: if a particle does not have a matching particle in the input PRTLoader object, it will be deleted.

# **Particle Force operator**

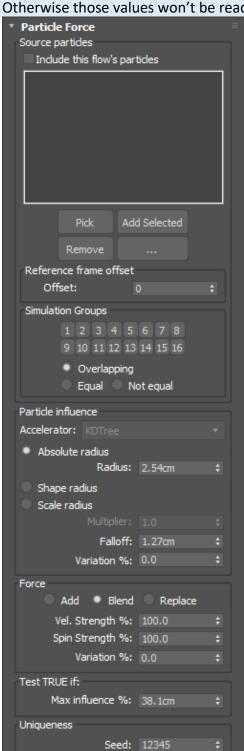


The Particle Force operator allows you to influence particle motion using particles from another flow.

### INFO:

A common application of this effect is to upsample simulations, by driving a high-resolution flow with the motion computed in a lower resolution flow.

**Note:** Please ensure you've enabled both spin and velocity caching in any referenced **tyFlow** objects! Otherwise those values won't be read properly by the operator and you will get incorrect result



**Source Particles** 

**Include this flow's particles**: the current flow's particles will contribute forces.

**Particle list**: the input particle objects which will be used to derive particle forces.

**Offset**: the frame offset that the input flow object will be evaluated at.

**Simulation groups**: controls which particle simulation groups will be read from the input flow objects.

### **Particle Influence**

**Accelerator**: the nearest-neighbor search algorithm used to find particles in the input flow that are closes to particles in the current event.

**Absolute radius**: the radius of each particle force will be set to a specific value.

**Radius**: the specific radius value.

**Shape radius**: the radius of each particle force will be set to each particle's shape mesh radius.

**Scale radius**: the radius of each particle force will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

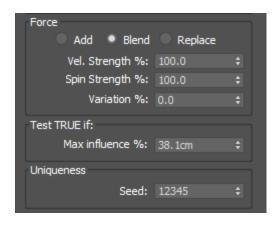
**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Variation %:** the per-particle percentage of variation to apply.

### TIP:

When upsampling a simulation using the Particle Force operator, a good rule of thumb is to set "radius" to roughly 75-100% of the radius of the input particles, and to set "falloff" to roughly 50% of the radius of the input particles.

### **Particle Force Continued**



### **Force**

**Add/Blend/Replace**: controls how forces computed by the operator will affect existing particle velocities.

**Vel strength %**: the percentage of the computed force to apply to particle velocity channels.

**Spin strength** %: the percentage of the computed force to apply to particle spin channels.

**Variation** %: the per-particle percentage of variation to apply.

## **Test TRUE If**

**Max influence**: particles will satisfy the TRUE condition if the combined max influence of input particles (based on the radius and falloff values) is less than this value.

## Uniqueness

# **Particle Groups operator**



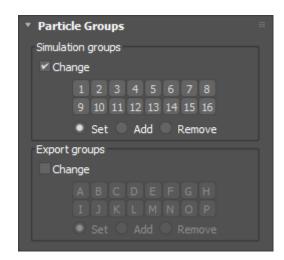
The Particle Groups operator can be used to set simulation/export groups on particles. The group settings can then be used to filter particles from simulation/export.

## INFO:

A particle group is a set of assignable bit flags. Two groups overlap if they share one or more of the same flags. Groups with no flags will overlap with other flagless groups, but if a group has no flags it will not overlap with a group that has one or more flags set. If two groups overlap, they pass the group filter test.

For example, a group set to [1] will overlap with a group set to [1][2]. But a group set to [1] will not overlap with a group set to [2].

Understanding this principle makes it very easy to filter particles later in a simulation, or in **tyFlow's** various export functions. For example, in a Particle Physics operator you could use groups to ensure that only certain particles will exert forces on each other. Groups can be utilized all throughout **tyFlow** and play an important role in overall flow organization.



### **Simulation groups**

Simulation groups can be utilized within simulations themselves, and have no effect on particles that are exported using the various export functions.

**Change**: controls whether this operator will modify existing particle simulation groups.

**Simulation groups**: controls which particle simulation groups will be assigned to particles.

**Set/Add/Remove**: controls which operation will be used to assign simulation groups.

## **Export groups**

Export groups can be utilized within particle export functions, and have no direct effect on simulations.

**Change**: controls whether this operator will modify existing particle export groups.

**Export groups**: controls which particle export groups will be assigned to particles.

**Set/Add/Remove**: controls which operation will be used to assign export groups.

# **Path Follow operator**



The Path Follow operator allows you to apply forces to particles using regular shapes and splines.

### TIP:

The default additive nature of forces can cause particles to quickly fly away from any input shapes. In order to cause particles to follow shapes more closely, force damping is required. The best way to dampen forces is to add a Slow operator to your event (above this operator) and increase the velocity slowdown value until the desired effect is achieved.

### **Path Follow Rollout Top Half**



### **Shapes**

**Shape object list**: the list of input shape objects.

### **Follow**

Follow forces are forces that are parallel to shape tangents.

**Velocity:** the value of the force applied to particles.

**Variation** %: the per-particle percentage of variation to apply.

**Falloff**: the distance over which forces will be diminished, according to the inverse-square law. A value of 0 means no falloff will be computed.

**Accel %**: controls how quickly particle forces will accelerate from their old values to the newly applied values.

### Attraction

Attraction forces are forces that are perpendicular to shape tangents, which cause particles to move toward shapes.

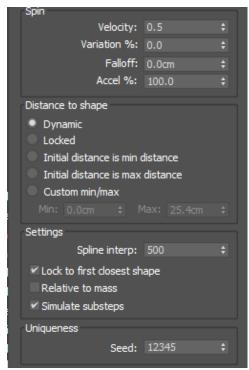
**Velocity:** the value of the force applied to particles.

**Variation** %: the per-particle percentage of variation to apply.

**Falloff**: the distance over which forces will be diminished, according to the inverse-square law. A value of 0 means no falloff will be computed.

**Accel** %: controls how quickly particle forces will accelerate from their old values to the newly applied values.

#### **Path Follow Rollout Bottom Half**



## Spin

Spin forces are forces that are perpendicular to shape tangents, which cause particles to move around shapes.

**Velocity:** the value of the force applied to particles.

**Variation** %: the per-particle percentage of variation to apply.

**Falloff**: the distance over which forces will be diminished, according to the inverse-square law. A value of 0 means no falloff will be computed.

**Accel** %: controls how quickly particle forces will accelerate from their old values to the newly applied values.

## Distance to shape

**Dynamic**: no constraints are placed on the distance particles are allowed to travel towards/away from input shapes.

**Locked** the initial distance any given particle is to an input shape, is the exact distance they must maintain while affected by the operator.

**Initial distance is min distance** the initial distance any given particle is to an input shape, is the minimum distance they must maintain while affected by the operator. Particles are allowed to travel further than that distance away from the input shapes.

**Initial distance is max distance** the initial distance any given particle is to an input shape, is the maximum distance they must maintain while affected by the operator. Particles are allowed to travel closer than that distance towards the input shapes.

Custom min/max: the range of distances a particle must stay within, from input shapes, is defined by the user.

Min/Max: the range of distances used in 'Custom min/max' mode"

### **Settings**

**Spline interp**: the number of sub-segments to slice input shapes up into, which will be used to accelerate the shape proximity algorithm.

**TIP:** If shapes have a lot of overall curve complexity, this value should be increased.

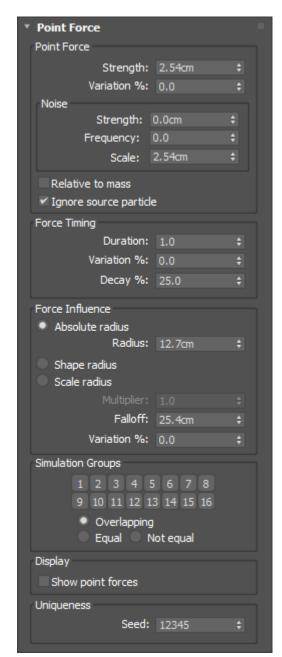
- Lock to first closest shape: controls whether particles will continually look for a new nearest shape to derive their forces from, or if they'll stay locked to the same shape for their entire stay within the event.
- **Relative to mass**: controls whether particle forces will be relative to particle masses.
- **Simulate substeps**: forces will be interpolated in a way that simulates the addition of forces at smaller simulation substeps.

### Uniqueness

# **Point Force operator**



The Point Force operator allows you to inject radial forces into a simulation, at particle locations.



### **Point Force**

**Strength:** the strength of each point force.

**Variation** %: the per-particle percentage of variation to apply.

### **Noise**

Allows you to add turbulence to the point forces.

**Strength**: the strength of the turbulent noise.

**Frequency**: the frequency of the turbulent noise.

Scale: the scale of the turbulent noise

**Relative to mass**: the force applied to a particle will be relative to its mass.

**Ignore source particle**: the particle from which the force originated will not be affected by the force.

### **Force Timing**

**Duration**: the duration, in frames, each point force will be active.

**Variation** %: the per-particle percentage of variation to apply.

**Decay**: the amount of decay to apply to forces, each time step.

**INFO:** Decay will decrease force strength over time.

### **Force Influence**

**Absolute radius**: the radius of each point force will be set to a specific value.

**Radius**: the specific radius value.

**Shape radius**: the radius of each point force will be set to each particle's shape mesh radius.

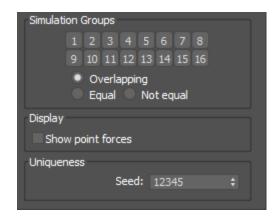
**Scale radius**: the radius of each point force will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Variation** %: the per-particle percentage of variation to apply.

#### **Simulation Groups**



**Simulation groups**: controls which particle simulation groups will be affected by the point forces.

## Display

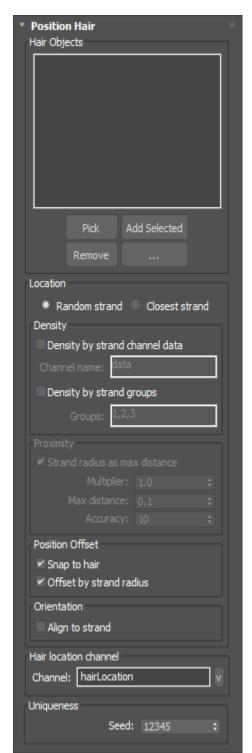
**Show point forces**: controls whether point forces will be visualized in the viewport, as spheres representing the size of their influence in the scene.

## Uniqueness

# **Position Hair operator**



The Position Hair operator allows you to position particles on Ornatrix hair strands



## **Hair Objects**

**Hair objects list**: the list of objects which have Ornatrix hair modifiers applied to them.

### Location

**Random strand**: particles will be uniformly placed on random Ornatrix hair strands.

**Closest strand**: particles will be placed on the closest Ornatrix hair strand, if they satisfy the distance condition.

**Note:** You can use the output functionality of this operator to send particles which satisfied the distance condition to another event.

### **Density**

These settings apply to "random strand" location mode.

**Density by strand channel data**: random strands will be weighted by Ornatrix channel data. The closer a vertex data value is to 1.0, the more likely it will receive a particle.

**Channel name**: the Ornatrix vertex data channel name.

**Density by strand groups**: only specified strand groups will receive particles.

**Groups**: the names of the Ornatrix strand groups, separated by commas.

### **Proximity**

These settings apply to "closest strand" location mode.

**Strand radius as max distance**: the maximum distance a particle can be to a strand before being placed on the strand is equal to the strand's radius.

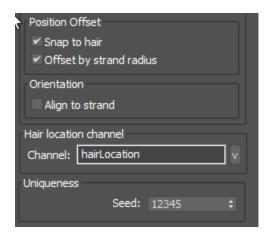
**Multiplier**: a multiplier to apply to the strand radius value.

**Max distance**: the maximum distance a particle can be to a strand before being placed on a strand is equal to this absolute value.

**Accuracy**: the maximum number of nearby hair strand vertices to examine when doing the hair strand proximity search.

### **INFO:**

The Position Hair operator uses an accelerated proximity search algorithm that first searches for nearest strand vertices, and from the list of those vertices then checks the distance between the particle and the strand segments adjacent to the nearest vertices. Setting the accuracy setting too low can result in particles being assigned to the wrong strand segments. It is recommended to keep the accuracy setting to between 10 and 25. Alternatively, increasing the segment count of the hairs themselves can improve search accuracy.



## **Position Offset**

**Snap to hair**: particles moved to strands will be snapped to the closest point on the strand.

**Offset by strand radius**: the minimum distance a particle can be from a strand is the radius of the strand.

## **Orientation**

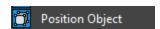
**Align to strand**: particles will be aligned to the strands they are placed on, with the strand tangent being the forward vector of the orientation matrix, and the vector to the closest point on the strand being the up vector of the orientation matrix.

### Hair location channel

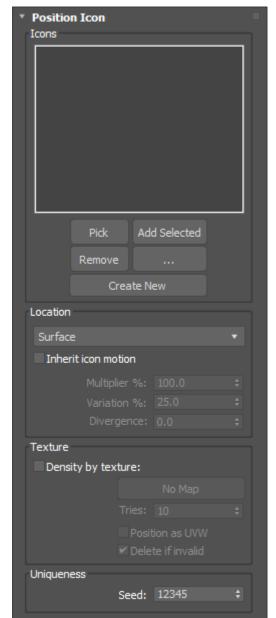
**Channel**: the custom TM channel where data required for the Hair Bind operator should be stored.

## Uniqueness

# **Position Icon operator**



The Position Icon operator allows you to position particles in the scene relative to a **tyIcon** object.



### **Icons**

**Icon object list**: the list of **tyIcon** objects which will be used to position particles in the scene.

### Location

**Location type**: the location on the **tyIcon** objects where particles will be placed.

**Inherit icon motion**: controls whether particles will inherit the velocity of their **tyIcon**.

**Multiplier** %: the amount of velocity particles will inherit.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: the angle of divergence applied to inherited velocity vectors.

### **Texture**

**Density by texture**: controls whether particle densities on a **tyIcon** will be controlled by a texture map.

**Texmap**: the texture map which will control particle densities.

**Tries**: the number of attempts to make, while searching for a valid location on the **tyIcon** (based on the texture map), before stopping the search.

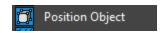
**Position as UVW**: the position of the particle in world-space will be used as the UVW value used to sample the density texture

**INFO:** If "position as UVW" is disabled, implicit UVWs will be assigned to **tylcons**, with corner-to-corner UVW values ranging from 0.0 to 1.0. If location type is set to "volume", the implicit UVWs will be 3-dimensional to account for the **tylcon's** depth.

**Delete if invalid**: controls whether particles that cannot find a valid location on the **tyIcon** are deleted.

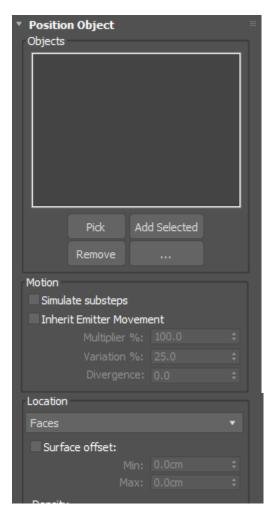
### Uniqueness

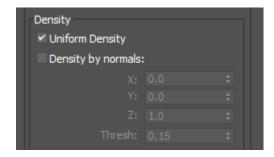
# **Position Object operator**



The Position Object operator allows you to position particles on the surface or in the volume of a scene object's geometry.

## **Position Object Rollout Top Section**





## **Objects**

**Object list**: the list of scene objects on whose geometry particles will be scattered.

### **Motion**

**Simulate substeps**: positions will be interpolated in a way that simulates the addition of positions at smaller simulation substeps.

**Inherit emitter movement**: controls whether particles will inherit velocities from the object on which they're scattered.

**Multiplier**: the amount of influence an object's velocity will have on particle velocities.

**Variation %**: the per-particle percentage of variation to apply.

**Divergence**: the angle of divergence applied to inherited velocity vectors.

### Location

**Location type**: the location on the scene object geometry where particles will be placed.

**Surface Offset**: controls whether particles will be offset from the surface they are placed on.

**Min/Max**: the minimum/maximum offset a particle may have from the surface it is placed on.

**Note:** Multiple density conditions can be enabled simultaneously to have an overall cumulative effect on where particles will be distributed.

**Uniform Density**: the number of particles that will be scattered on a particular surface mesh triangle will be relative to that triangle's total area, resulting in an overall uniform distribution of particles on the surface mesh.

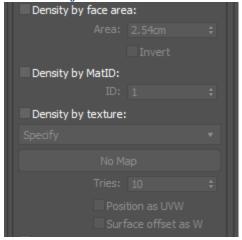
### **Density By Normals**

**Density by normals**: controls whether the normal direction of surface mesh triangles will affect particle distribution.

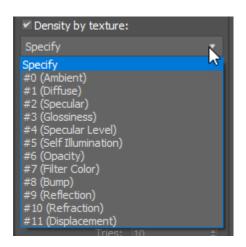
**XYZ**: The target normal direction vector.

**Thresh**: the threshold for particle distribution, given a particular surface mesh triangle's face normal in comparison to the target vector.

#### **Position Object Rollout Middle Section**



#### **Density by Texture Selection Box Detail**



## **Density By Face Area**

**Density by face area**: controls whether surface mesh triangle areas will affect particle distribution.

**Area**: the minimum face area required for a particular surface mesh triangle to quality for particle distribution.

**Invert**: the minimum area threshold will instead be treated as a maximum area threshold.

## **Density By MatID**

**Density by matID**: controls whether surface mesh triangle material IDs will affect particle distribution.

**ID**: the material ID value that a surface mesh triangle must have in order to quality for particle distribution.

### **Density By Texture**

**Density by texture**: controls whether a texture sampled on the surface of an input object will affect particle distribution.

**Texmap source**: the source of the texture map. If set to "specify", it will be the specified texture map below, otherwise it will be derived from the selected channel of the position object's material.

**Texmap**: the specified texture map which will control particle densities.

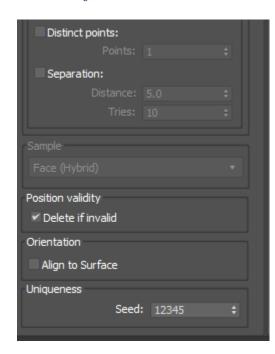
**Tries**: the number of attempts to make, while searching for a valid location on an input mesh (based on the texture map), before stopping the search.

**Position as UVW**: the position of the particle in world-space will be used as the UVW value used to sample the density texture.

**Surface offset as W**: the distance from the particle to the closest surface will be used as the W coordinate in the UVW value which is used to sample the texture.

**Note:** "Surface offset as W" can be beneficial for sampling 3D maps like noise maps, but it can be detrimental for 2D maps like the Vertex Color map, which interprets any value over 1 in any UVW coordinate as solid white. "Surface offset as W" should be disabled when using Vertex Color maps for particle density.

#### **Position Object Rollout Bottom Section**



#### **Distinct Points**

**Distinct points**: controls whether particle locations will be limited to a finite number of locations on the input objects.

**Points**: the number of distinct points to use.

## **Separation**

**Separation**: controls whether particles must remain a certain distance from other particles before their location on a surface is considered valid.

**Distance**: the minimum distance allowed between particles on a surface.

**Tries** the number of attempts to make, while searching for a valid location on an input mesh (based on the separation distance), before stopping the search.

## Sample

**Sample type**: controls which sampler will be used to determine closest-object proximities for particles.

## **Position Validity**

If particle exceeds the max number of tries allowed for a particular density condition, its resulting position will be considered invalid.

**Delete if invalid**: controls whether particles with invalid positions will be deleted.

## Orientation

**Align to surface**: aligns particles to the surface of their input object.

## TIP:

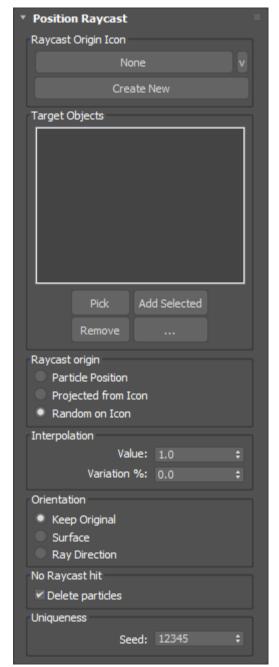
The Position Object operator offers no extra controls over surface alignment. A Rotation operator should be used instead if more control over surface alignment is required.

# Uniqueness

# **Position Raycast operator**



The Position Raycast operator allows you to cast rays towards the surface of an object, and place particles where the rays hit.



### **Raycast Origin Icon**

**tyIcon object**: the **tyIcon** object used to cast the rays. Rays will be cast in the direction of the **tyIcon's** arrow.

## **Target Objects**

**Target object list**: the list of input objects whose surface meshes will be hit by rays.

## Raycast origin

**Particle Position**: rays will be cast from the particle's current position.

**Project from Icon**: rays will be cast from the particle's current position, projected onto the plane of the applicable tyIcon.

**Random on Icon**: rays will be cast from a random location on the enclosed plane of the applicable tylcon.

## Interpolation

**Value**: the amount to interpolate particle positionx from their previous value to the new value.

**Variation %:** the per-particle percentage of variation to apply.

#### Orientation

**Keep original**: particle orientations will not be changed.

**Surface**: particles will be aligned to the surface their ray hit.

#### TIP:

The Position Raycast operator offers no extra controls over surface alignment. A Rotation operator should be used instead if more control over surface alignment is required.

**Ray direction** particles will be aligned to the direction of the rays.

## No Raycast Hit

**Delete Particles**: if enabled, particles whose rays do not hit an object will be deleted.

#### Uniqueness

# **Position Transfer operator**



The Position Transfer operator allows you to transfer particles from one surface to the exact relative location on another surface with the same topology.

#### TIP:

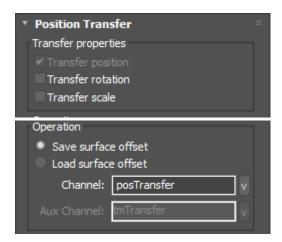
The Position Transfer operator can be especially helpful when you want to simulate particles on a static surface, and then transfer their motion to a matching animated surface (ie, an animated character). It is often easier to perform this kind of transfer from static to animated mesh, than to try and perform the initial particle simulation on the animated mesh directly.

#### INFO:

For the position transfer to work correctly, the objects in the list must have the same topology and be in the same order in all matching Position Transfer operators.

#### NOTE:

When transferring position data between flows, don't forget to enable Custom TM channel caching in the source flow.



## **Transfer properties**

**Transfer position/rotation/scale**: controls which relative particle properties to transfer between surfaces.

## **Operation**

**Save surface location**: the operator will save the relative surface location of particles to a custom TM data channel.

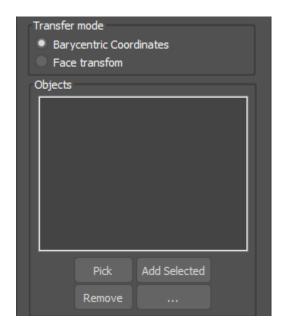
**Load surface location**: the operator will load the relative surface location of particles from a custom TM data channel.

**Channel**: the main custom TM data channel where transfer data will be stored.

**Aux channel**: the secondary custom TM data channel where rotation/scale offset data will be stored, if rotation/scale property transfer is enabled.

## NOTE:

If both channels are required for the transfer (ie, if you've enabled rotation/scale transfer), ensure they are not assigned the same name.



#### **Transfer Mode**

**Barycentric Coordinates**: saves the transfer data as relative barycentric coordinates of the closest face on the source mesh.

**Face transform**: saves the transfer data as a position relative to a transform constructed from the closest face on the source mesh.

#### INFO:

"Barycentric coordinate" mode works best if particles are near the source surface and the source surface is fairly smooth. It does a good job of smoothly interpolating particles between faces on the target mesh, so long as particles are not too far away from the source surface and the surface does not have a lot of hard edges/angles, otherwise it can result in popping artifacts. "Face transform" mode will not smoothly interpolate particles between faces, but does a better job at storing the absolute particle position relative to the source mesh, even if the particles are heavily offset from the source surface.

#### NOTE:

When transferring rotation/scale data, face transform mode must be used.

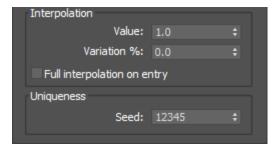
#### **Objects**

**Object list**: the list of input objects that particles will be transferred to/from.



#### Sample

**Sample type**: controls which sampler will be used to determine closest-surface proximities for particles.



## Interpolation

**Value**: the amount to interpolate particle positions from their previous value to the new value.

**Variation %:** the per-particle percentage of variation to apply.

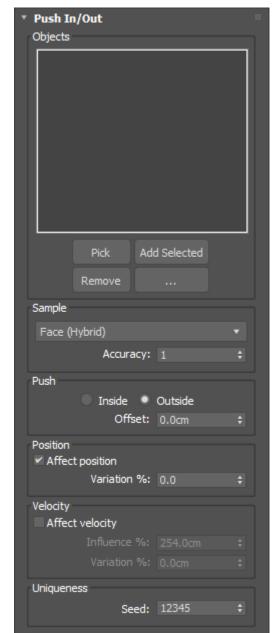
**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

## Uniqueness

# Push In/Out operator



The Push In/Out operator allows you to push particles in/out of closed surfaces.



#### **Objects**

**Object list**: the list of objects whose volumes will be used for in/out tests.

### **Sample**

**Sample type**: controls which sampler will be used for in/out tests.

**Accuracy**: controls the accuracy of the raycaster used to compute information about whether a particle is inside or outside of an object's volume.

#### **Push**

**Inside/Outside**: the push operation to perform. "Inside" will push particles inside a surface, whereas "outside" will push particles outside a surface.

**Offset**: the amount of additional offset to apply to target particle locations, along the direction they are moved in/out of a surface.

**Note:** When offset is set to value other than 0, the operator will begin to also affect particles within that distance range to the surface. For example, if push mode is set to "outside" and offset is set to "10", any particles within the surface or within a distance of 10 units to its surface will be affected.

#### **Position**

**Affect position**: controls whether particles will be forcefully moved in/out of volumes.

**Variation** %: the per-particle percentage of variation to apply.

## Velocity

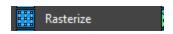
**Affect velocity**: controls whether particle velocities will modified with the vector required to move them in/out of a surface.

**Influence**: the amount of influence the in/out vector will have on a particle's velocity.

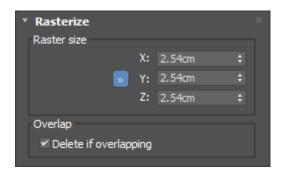
**Variation %:** the per-particle percentage of variation to apply.

## Uniqueness

# **Rasterize operator**



The Rasterize operator allows you to clamp particle positions to the cells of an implicit grid. This is similar to what the Birth Voxels operator does, except this doesn't birth any new particles – instead, it changes the locations of existing particles.



## **Raster Size**

**X/Y/Z**: the size of the cells of the implicit grid.

## Overlap

**Delete if overlapping**: controls whether multiple particles are allows in the same cell. If enabled, only 1 particle will be allowed in a cell – the rest will be deleted.

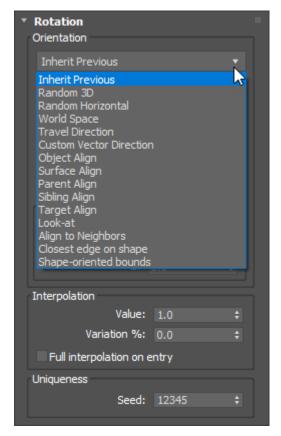
# **Rotation operator**



The Rotation operator allows you to control particle orientations.

## Rotation Rollout – Orientation Types

**Rotation Rollout Selection Drop Down Choices** 



**Orientation type**: the starting orientation value, prior to any further modifications.

**Inherit previous**: uses the current value of a particle's orientation as the start value.

**Random 3D**: uses a random orientation as the start value.

**Random horizontal**: uses a random orientation around the world's Z-Axis as the start value.

World space: uses a world-aligned orientation as the start value.

**Travel direction**: uses an orientation aligned to the particle's velocity as the start value.

**Custom vector direction**: uses an orientation aligned to a custom data channel as the start value.

**Surface align**: uses an orientation aligned to the closest point on a scene object's surface mesh as the start value.

**Parent align**: uses an orientation aligned to the parent of each particle, if the particle has a parent.

**Sibling align**: uses an orientation aligned to the adjacent sibling of each particle, if such a sibling exists.

**Look-at**: uses an orientation directed towards the closest point on an input object as the start value.

**Align to neighbors**: uses an orientation directed towards the particle's neighbors as the start value.

**Closest edge on shape**: the orientation will be derived from the vector defined by the closest edge on the particle shape to the particle location.

**Shape-oriented bounds**: the orientation will be derived from the oriented bounding box of the particle shape mesh. The longest axis of the bounding box will be the up vector of the rotation matrix, and the second longest axis will be the forward vector.

## **Rotation Rollout**



**Affect particle shape orientation**: controls whether a particle's shape mesh orientation will be affected by the particle's orientation change.

#### INFO:

Disabling this setting has the effect of orientating the particle's transform without changing the orientation of the particle's mesh. In other words, it's akin to rotating the particle's pivot.

**Note:** Disabling "affect particle shape orientation" can have a big impact on memory usage and performance, because each particle's mesh will have to be manually adjusted in order to compensate for the particle's orientation change.

**X/Y/Z**: controls the amount of offset applied to starting orientations, along each axis.

**Divergence**: controls the degrees of random divergence to apply to starting orientations.

**Restrict diverg. to axis**: restricts the amount of divergence applied to each axis.

## **Divergence Axis:**

**X/Y/Z**: the amount of divergence to apply to each axis.

#### **Interpolation**

**Value**: the amount to interpolate particle orientations from their previous value to the new value.

**Variation** %: the per-particle percentage of variation to apply.

**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

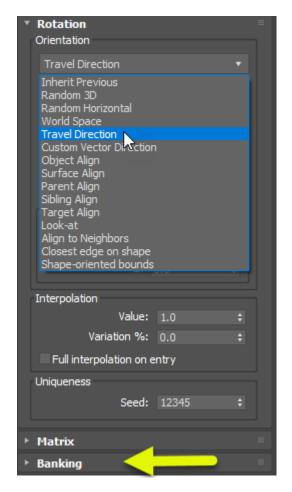
#### TIP:

In order to animate particles rotating towards a particular target orientation, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.

## Uniqueness

# **Banking Rollout**

The "Banking Rollout" appears when rotation type is set to **Travel Direction** 





# Banking

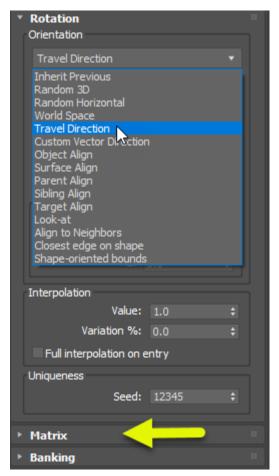
**Multiplier**: the amount of banking to apply to starting orientations, around the particle's travel direction.

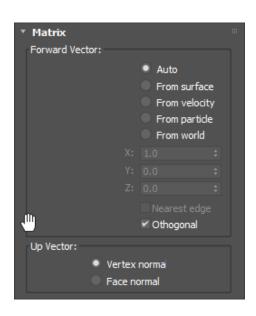
**Smooth**: the amount of temporal smoothing to apply to banking values, based on changes to the particle's travel direction over time.

## **Matrix Rollout**

The "Matrix Rollout" appears when rotation type is set to either of these modes:

## Travel Direction / Custom Vector Direction / Surface Align / Look-At / Closest Edge on Shape





**Matrix Rollout** 

These settings control how secondary vectors of the orientation matrix will be derived (the up vector is the main alignment vector, and the right vector is derived from the cross product between the up and forward vectors).

#### Forward Vector:

**Auto**: the forward vector of the surface alignment transforms will be automatically computed based on the orientation of the up vector.

**From surface**: the forward vector of surface alignment transforms will be derived from the surface itself.

**From velocity**: the forward vector of surface alignment transforms will be derived from each particle's velocity direction.

**From particle**: the forward vector of surface alignment transforms will be derived from each particle's transform forward vector.

**From world**: the forward vector of surface alignment transforms will be defined in world space.

**X/Y/Z**: allows users to define world-space forward vectors, when "from world" is enabled.

**Nearest edge**: when forward vector is derived from surface, the forward vector will be set to the nearest edge on the nearest face to the particle position, rather than the first edge.

**Orthogonal**: controls whether surface alignment orientations will be made orthogonal to the up vector.

## **Up Vector:**

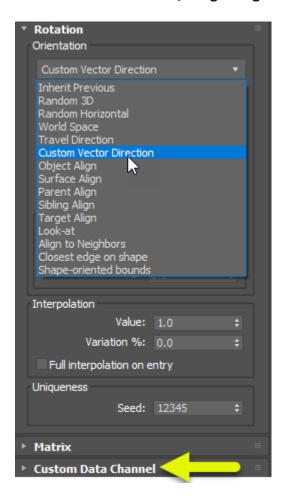
**Vertex normal**: controls whether the up vector of surface alignment transforms will be derived from the nearest vertex normal.

**Face normal**: controls whether the up vector of surface alignment transforms will be derived from the nearest face normal.

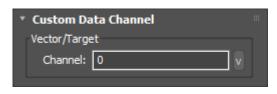
## **Custom Data Channel Rollout**

The "Custom Data Channel Rollout" appears when rotation type is set to either of these modes:

## **Custom Vector Direction / Target Align**



These settings pertain to rotation types which rely on the input of a custom data channel.

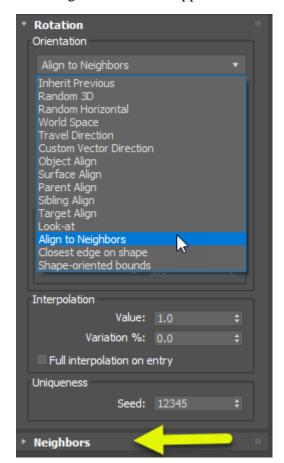


## Vector/Target

**Channel**: the channel to derive the custom vector direction from.

# **Neighbors Rollout**

The "Neighbors Rollout" appears when rotation type is set to Align to Neighbors



These settings control neighbor searches in neighbor align mode.



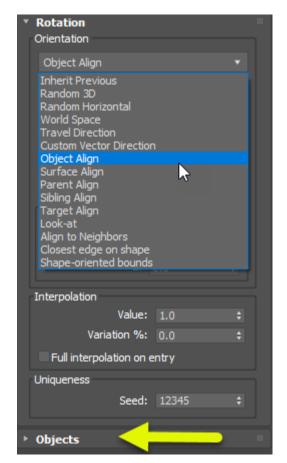
**Radius**: The particle neighbor search radius. Nearby particles within this radius will be considered neighbors.

**Simulation groups**: controls which particle simulation groups nearby particles must match in order to be considered neighbors.

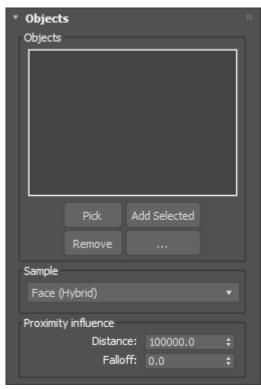
# **Objects Rollout**

The "Objects Rollout" appears when rotation type is set to either of these modes:

## Object Align / Surface Align / Look-At



These settings control which objects will be sampled for various rotation types, and the proximity of their influence.



## **Objects**

**Object list**: the list of input objects used for surface alignment or facing mode.

#### **Sample**

**Sample**: controls which sampler will be used for surface alignment.

## **Proximity influence**

The proximity influence extends outwards from nearby objects

**Distance**: particles within this distance will be fully affected.

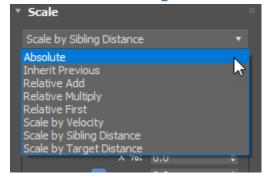
**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

# **Scale operator**



The Scale operator allows you to control the size of particles.

# Scale Rollout – Scaling Methods



**Scale type**: the particle scaling method to use.

**Absolute**: particle scales will be set to a user-defined absolute value.

**Inherit previous**: uses the current value of a particle's scale as the starting point, prior to modifications.

**Relative add**: add a user-defined scale value to the current value of the particle's scale.

**Relative multiply**: multiplies the current value of the particle's scale by a user-defined vector.

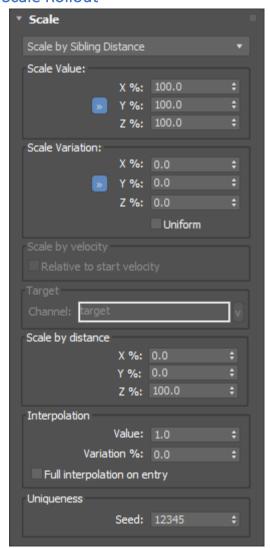
**Relative first**: multiplies the value of the particle's scale when it first entered the operator by a user-defined vector.

**Scale by velocity**: constructs a scale vector for each particle using its velocity as a multiplier along each axis. The multiplier is applied to the scale of the particle when it enters the event. Scale X/Y/Z values set to 0 will maintain the original scale of each particle.

**Scale by Sibling Distance**: the z-axis of the particle's scale vector will be set to the distance between the particle and its adjacent sibling.

**Scale by Target Distance**: the z-axis of the particle's scale vector will be set to the distance between the particle and its target.

## Scale Rollout



#### Scale value

**X/Y/Z** %: the absolute values to use for the scale modification vector.

#### INFO:

The effect of the absolute values on a particle's scale depends on which scale type is selected.

#### **Scale variation**

**X/Y/Z** %: the amount of variation to apply along each axis of the intermediate scale values.

**Uniform**: the variation applied along each axis will be uniform, preventing proportion skewing.

## Scale by velocity

**Relative to start velocity**: the velocity-scale multiplier will be relative to the velocity of the particle when it entered the event, rather than zero.

## Scale by distance

**X/Y/Z**: controls the level of influence the scale-by-distance modes will have on each axis of a particle's scale vector.

#### Interpolation

**Value**: the amount to interpolate particle scales from their previous value to the new value.

**Variation %**: the per-particle percentage of variation to apply.

**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

## TIP:

In order to animate particles scaling towards a particular target scale, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.

# Uniqueness

# **Texmap Rollout**

The Texmap Rollout allows you to control particle scales by sampling textures on nearby surfaces and multiplying scales by the resulting values. The darker the sampled texture value, the smaller the scale multiplier.



**Multiply by texmap**: controls whether texmap scaling will be enabled.

**Texmap**: the texture map to sample.

**Surface offset as W**: the distance from the particle to the closest surface will be used as the W coordinate in the UVW value which is used to sample the texture.

**Object list**: the list of objects whose surfaces will be sampled for UVW coordinates.

**Sample**: controls which sampler will be used for surface proximity tests.

**Affect X/Y/Z**: provides control over the amount of influence a texmap will have on each axis of the resulting scale vector.

# **Proximity Rollout**

The proximity rollout allows you to control particle scales by measuring distances to nearby scene objects and their surfaces, and multiplying scales by the resulting values.



**Multiply by proximity**: controls whether proximity scaling will be enabled.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**Invert**: inverts the affect of the proximity scale multiplier.

**Object list**: the scene objects used to calculate proximity multipliers.

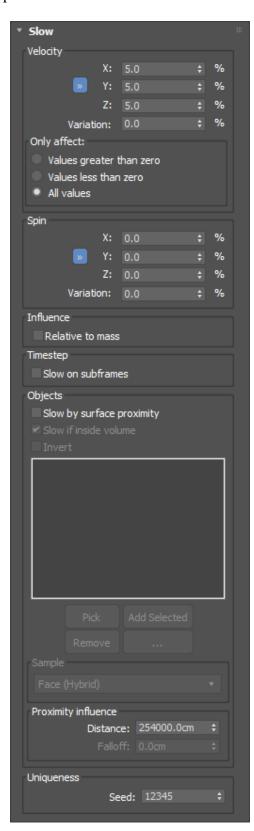
**Sample**: controls which sampler will be used for surface proximity tests.

**Affect X/Y/Z**: provides control over the amount of influence the proximity multiplier will have on each axis of the resulting scale vector.

# **Slow operator**



The Slow operator allows you to dampen particle forces. It should be used instead of a Drag spacewarp to slow particles down within a flow.



#### TIP:

The Slow operator should usually be placed *above* any other operators that apply forces to particles within an event. Since operators are evaluated in order, this will dampen forces when the next time step is evaluated, without damping forces added to particles in the current time step.

### Velocity

**X/Y/Z**: the amount to dampen particle velocities, per axis.

**Variation**: the per-particle amount of variation to apply.

#### **Only Affect:**

**Values greater than zero**: a velocity value along a particular axis will only be dampened if it is above zero.

**Values less than zero**: a velocity value along a particular axis will only be dampened if it is below zero.

All values: all velocities values will be dampened.

#### TIP:

Sometimes you might only want to dampen particle velocities in a certain direction. For example, you might want to dampen particle velocities as particles move upwards, but not as they move downwards with gravity (to prevent unwanted bouncing). Use the appropriate "only affect" type to precisely control which values will be dampened.

#### **Spin**

**X/Y/Z**: the amount to dampen particle spin, per axis.

**Variation**: the per-particle amount of variation to apply.

#### **Influence**

**Relative to mass:** the amount of damping applied will be relative to a particle's mass.

### **Timestep**

**Slow on subframes::** controls whether to apply damping to particles on subframes. Turning this on will affect how quickly particles will be slowed, if simulation time steps are set to smaller than 1 frame.

#### **Slow Operator Rollout Continued**



#### **Objects**

**Slow by surface proximity**: controls whether the affect of the operator will be relative to a particle's proximity to nearby objects.

**Slow if inside volume**: controls whether the proximity effect will be applied to particles inside a surface.

**Invert** inverts the effect of proximity slowing.

**Object list**: the list of input objects whose proximity will be tested.

# **Sample**

**Sample**: controls which sampler will be used for surface proximity tests.

## **Proximity influence**

The proximity influence extends outwards from the surface of nearby obstacles.

**Distance**: particles within this distance will be fully affected.

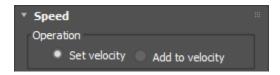
**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

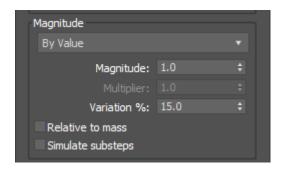
## Uniqueness

# **Speed operator**



The Speed operator allows you to assign velocities to particles. Velocity vectors are constructed by multiplying a direction vector by a magnitude.





## **Operation**

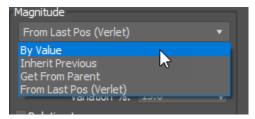
**Set velocity**: changes particle velocity to the value created by the Speed operator.

**Add to velocity**: adds the value created by the Speed operator to the existing particle velocity value.

#### Magnitude

**Velocity magnitude type**: the method used to define the desired magnitude of the resulting velocity value.

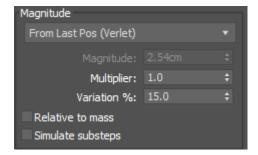
By Value: the magnitude will be set to an absolute value.



**Inherit Previous**: the magnitude will be set to the particle's current velocity magnitude.

**Get From Parent**: the magnitude be set to magnitude of the particle's parent particle velocity, if it has a valid parent.

**From Last Pos (Verlet)**: the magnitude will be the distance between the particle's current position and its position at the previous time step.



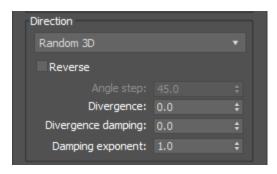
**Magnitude**: the absolute magnitude value.

**Variation** %: the per-particle percentage of variation to apply.

**Relative to mass**: controls whether magnitude values will be relative to a particle's mass.

**Simulate substeps**: forces will be interpolated in a way that simulates the addition of forces at smaller simulation substeps.

#### **Speed Operator Rollout Continued**



## **Direction Drop Down Menu Choices**



#### NOTE:

Divergence damping is the tendency for the magnitude of a speed vector to approach zero, as it diverges from its original direction. Increasing this value will slow particles down, depending on how much their direction vector diverges.

#### Direction

**Velocity direction vector type**: the method used to define the desired direction vector of the resulting velocity value.

**Along Icon Arrow**: the direction vector will be aligned to the nearest **tyIcon** object's arrow.

**Icon Center Out**: the direction vector will be aligned to a vector pointing outward from the center of the nearest **tyIcon** object.

**Icon Axis Out**: the direction vector will be aligned to a vector pointing outward from the central axis of the nearest **tyIcon** object.

**Random 3D**: the direction vector will be a random 3D vector.

**Random Horizontal**: the direction vector will be a random 3D vector around the world's Z-Axis.

**Inherit Previous**: the direction vector will be set to the particle's current velocity direction vector.

**Particle X/Y/Z**: the direction vector will be set to a local axis of the particle's orientation.

**Get From Parent**: the direction vector will be set to direction vector of the particle's parent particle velocity vector, if it has a valid parent.

**Outwards From Parent**: the direction vector will be set to direction from the parent particle to the particle, if it has a valid parent.

**From Last Pos** (Verlet): the direction vector will be set to the direction between the particle's current position and its position at the previous time step.

**Discretize**: the direction vector will be set to a discretized vector derived from the particle's current direction vector.

**Absolute**: the direction vector will be set to the dominant axis of the particle's current direction vector.

**Reverse**: reverses the resulting direction vector.

**Angle step**: the maximum angle that a discretized vector may deviate from a world-aligned axis.

**Divergence**: the degrees of random divergence to apply to the resulting direction vector.

**Divergence damping**: This value controls how much divergence damping to apply to resulting speed vectors.

**Damping exponent**: The dot product used to calculate the divergence damping value will be raised to this exponent. Increasing this value will increase damping for smaller divergences.

#### **Speed Operator Rollout Bottom Section**



#### **Noise**

Applies a turbulent noise multiplier to velocity values, based on particle positions in space.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

Scale: the scale of the noise.

Roughness: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

## Interpolation

**Value**: the amount to interpolate particle velocities from their previous value to the new value.

Variation %: the per-particle percentage of variation to apply.

#### **Icons**

**tyIcon object list**: the list of **tyIcon** objects to use for various values calculations. Only the closest object to each particle will be used.

## Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

# Spin operator



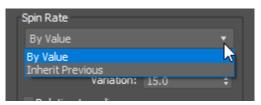
The Spin operator allows you to add spin motion to particles.



#### **Spin Rate**

**Spin magnitude type**: the method used to derive the magnitude of the spin vector.

By Value: the magnitude will be set to an absolute value.



**Inherit Previous**: the magnitude will be set to the particle's current spin magnitude.

**Spin rate**: the rate, in degrees per second, to spin particles.

**Variation**: the per-particle amount of variation to apply.

**Relative to radius**: the spin magnitude will be relative to a particle's radius.

**Radius**: the threshold radius value.

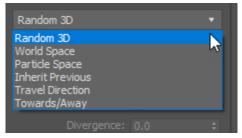
**Relative to velocity**: the spin magnitude will be relative to a particle's velocity.

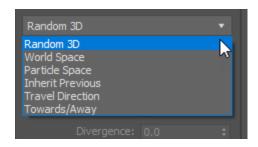
**Velocity**: the threshold velocity value.

**Relative to mass**: the spin magnitude will be relative to a particle's mass.

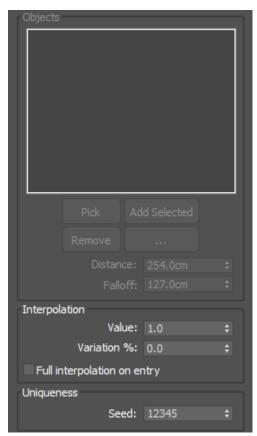
## **Spin Axis**

**Spin axis type**: the method used to derive the axis of the spin vector.









Random 3D: the spin axis will be a random 3D vector.

**World Space**: the spin axis will be a user-defined vector in world-space coordinate.

**Inherit Previous**: the spin axis will be set to the particle's current spin axis.

**Travel direction**: the spin axis will be aligned to the particle's travel direction.

**Towards/Away**: the spin axis will be aligned towards/away the closest input object, depending on the sign of the magnitude.

**X/Y/Z**: the values of the user-defined spin axis vector.

**Allow mirror**: controls whether some particles will spin in the opposite direction.

**Divergence**: the degrees of random divergence to apply to the resulting axis vector.

## **Objects**

**Input object list**: the list of scene objects to use for towards/away spinning.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

### **Interpolation**

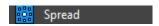
**Value**: the amount to interpolate particle spin values from their previous value to the new value.

**Variation** %: the per-particle percentage of variation to apply.

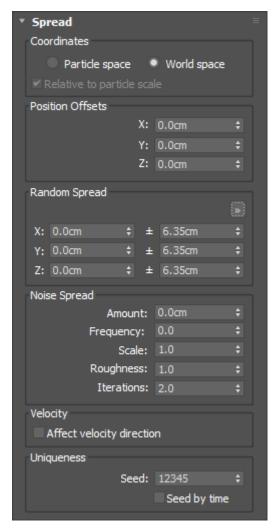
**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

## Uniqueness

# **Spread operator**



The Spread operator allows you to directly move particles in various directions, by adjusting their position instead of applying forces.



**Particle/World space**: controls which coordinate system the spread values will be calculated within.

**Relative to particle scale**: when in particle space mode, the spread values will be transformed relative to the particle's scale.

#### INFO:

When "relative to particle scale" is enabled, the scale of each axis of the particle's transform will act as a multiplier on the resulting spread vector. So if you offset a particle 1 unit along the z-axis, and the particle has a scale of [5,10,20], the particle will be moved 20 units up along the direction of its transform's z-axis in world space, even though in local particle space it will have only moved 1 unit along that axis.

**Position Offsets** 

**X/Y/Z**: allows you to move particles along a world-aligned axis by a precise value.

## **Random Spread**

**X/Y/Z**: the overall distance to move particles, in random directions.

 $\pm$  **X/Y/Z**: controls the amount of variation to apply to the initial X/Y/Z distances.

#### **Noise Spread**

The noise settings allow you to offset the way in which particle positions are moved.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

**Scale**: the scale of the noise.

**Roughness**: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

#### Velocity

**Affect velocity direction**: controls whether the direction between a particle's old position and its new position will affect the direction of its velocity vector.

## Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

# **Stop operator**



The Stop operator allows you to stop a particle from moving or spinning.

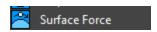


## **Stop Motion**

**Velocity X/Y/Z**: controls which axis of the velocity channel will be set to zero.

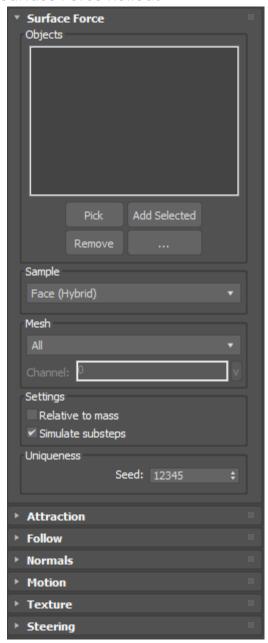
**Spin X/Y/Z**: controls which axis of the spin channel will be set to zero.

# **Surface Force operator**



The Surface Force operator allows you to add forces to particles using the surface properties of scene geometry.

## Surface Force Rollout



## Surface Force Rollout

**Object list**: the list of geometry input objects.

**Sample type**: controls which sampler will be used to determine closest-object proximities for particles.

**Mesh selection**: controls which meshes that forces will be derived from.

All: the closest mesh from the list will be used. **Random**: a random mesh from the list will be used. **Custom Float as Index**: a custom channel value will be used as a list index to determine which object's mesh to use.

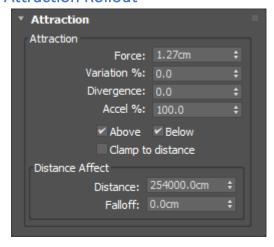
**Channel**: the custom float data channel to use.

**Relative to mass**: the amount of force applied will be relative to a particle's mass.

**Simulate substeps**: forces will be interpolated in a way that simulates the addition of forces at smaller simulation substeps.

#### Uniqueness

## **Attraction Rollout**



Attraction forces are forces which are directed from the particle location to the closest point on the target surface.

Force:: the amount of attraction force.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: the degrees of random divergence to apply to the resulting force.

**Accel %:** controls how quickly particle forces will accelerate from their old values to the newly applied values.

**Above/Below**: controls whether forces will be applied above/below surface normals.

**Clamp to Distance**: if the distance between a particle and the surface is less than the magnitude of the attraction force, the attraction force magnitude will be clamped to that distance.

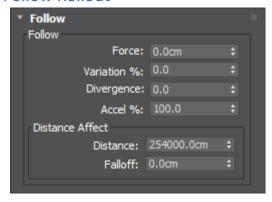
#### Distance affect

The distance affect extends outwards from the affecting surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

## **Follow Rollout**



Follow forces are forces which are parallel to the target surface.

**Force:**: the amount of follow force.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: the degrees of random divergence to apply to the resulting force.

**Accel** %: controls how quickly particle forces will accelerate from their old values to the newly applied values.

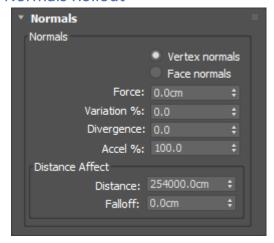
#### **Distance affect**

The distance affect extends outwards from the affecting surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

## **Normals Rollout**



Normal forces are forces which extend along the normal of the closest point on the target surface.

Force: the amount of normal force.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: the degrees of random divergence to apply to the resulting force.

**Accel** %: controls how quickly particle forces will accelerate from their old values to the newly applied values.

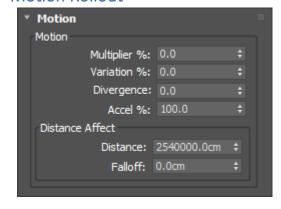
#### Distance affect

The distance affect extends outwards from the affecting surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law

## **Motion Rollout**



Motion forces are forces which are derived from the nearest surface velocity of the target surface.

Force: the amount of motion force.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: the degrees of random divergence to apply to the resulting force.

**Accel** %: controls how quickly particle forces will accelerate from their old values to the newly applied values.

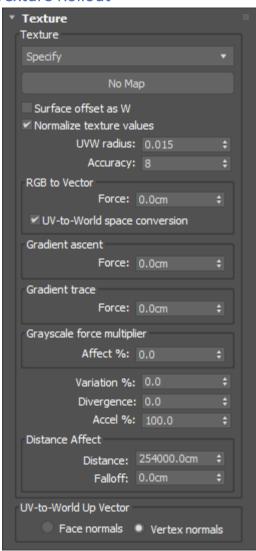
## Distance affect

The distance affect extends outwards from the affecting surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

## **Texture Rollout**



Texture forces are forces which are derived from texture map values sampled from the closest point on the target surface.

**Texmap**: the texmap to sample colors from.

**Surface offset as W**: the distance from the particle to the closest surface will be used as the W coordinate in the UVW value which is used to sample the texture.

**UVW Radius**: the samples radius (in UVW coordinates) to derive the gradient from.

**Accuracy**: the number of points to sampled in order to derive the gradient.

### **RGB** to Vector

RGB values sampled at the closest point on the surface to the particle will be used a world-space force vector.

**Force**: the amount of RGB force.

**UV-to-World space conversion**: when enabled, converts RGB values from UV space to world space. Disable this setting if RGB values already represent world-space force vectors.

#### **Gradient ascent**

Monochrome intensity values sampled around the closest point on the surface to the particle will be used to derive a directional gradient, transformed from UVW-space into world-space as a force vector.

#### TIP:

The direction of a gradient ascent force is from darker parts of a texture map to lighter parts of a texture map. If the monochrome intensity values in a particular area of the texture map are perfectly uniform, no force will be derived from that area.

**Force**: the amount of ascent force.

#### **Gradient trace**

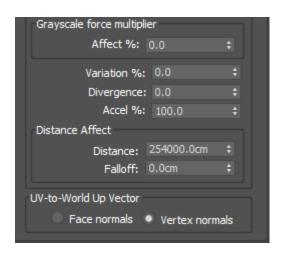
Monochrome intensity values sampled around the closest point on the surface to the particle will be used to derive a directional gradient. A perpendicular vector to that gradient will then be transformed from UVW-space into world-space as a force vector.

#### TIP:

The direction of a gradient trace force is around the borders of light/dark features of a texture map. If the monochrome intensity values in a particular area of the texture map are perfectly uniform, no force will be derived from that area.

**Force**: the amount of trace force.

#### **Texture Rollout Continued**



#### **Grayscale force multiplier**

Monochrome intensity values sampled around the closest point on the surface to the particle will be used to dampen accumulated Surface Force velocities. For example, a velocity of [1,0,0] will be changed to [0.75, 0, 0] if the sampled monochrome value is 0.75 and "affect" is set to 100%.

**Affect %:** the amount of influence that the grayscale force multiplier will have on accumulated Surface Force velocities.

**Variation** %: the per-particle percentage of variation to apply.

**Divergence**: the degrees of random divergence to apply to the resulting force.

**Accel %:** controls how quickly particle forces will accelerate from their old values to the newly applied values.

#### **Distance affect**

The distance affect extends outwards from the affecting surface.

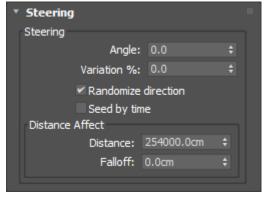
**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

## **UV-to-World Up Vector**

**Face/Vertex normals**: controls which surface normals will be used to compose the UV-to-World conversion matrix. Face normals are better for segmented/flat surfaces, Vertex normals are better for smooth/curved surfaces.

# **Steering Rollout**



Steering forces are forces which steering particles towards the cross product between the particle's velocity and the nearest surface normal.

**Angle**: the amount of steering force to apply to a particle's direction of travel.

**Variation** %: the per-particle percentage of variation to apply.

**Randomize direction**: controls whether the steering force can be on either side of the particle's velocity vector, relative to the nearest surface normal.

**Seed by time**: steering force variation will be reseeded at each time step.

#### Distance affect

The distance affect extends outwards from the affecting surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

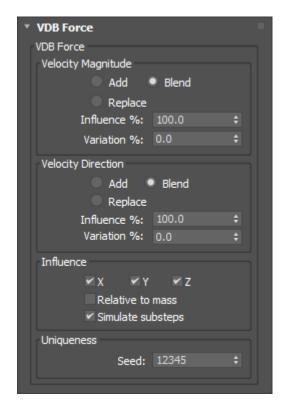
# **VDB** Force operator



The VDB Force operator allows you to apply forces from VDB voxels to particles.

### **INFO:**

This operator derives forces from VDB data directly inside the flow. If you want to use forces from a VDB setup in another flow, for the time being you will have to export the VDB data of that other flow to a .vdb file and bring it into the current flow using a supported container (PhoenixFD or FumeFX container) and a Fluid Force operator.



#### **Velocity Magnitude**

These controls affect the overall magnitude of VDB forces applied to particle velocities.

**Add/Blend/Replace**: controls how forces computed by the operator will affect existing particle velocities.

**Influence** %: controls how much influence VDB forces will have on particle velocities.

**Variation** %: the per-particle percentage of variation to apply.

## **Velocity Direction**

These controls affect the overall direction of VDB forces applied to particle velocities.

**Add/Blend/Replace**: controls how forces computed by the operator will affect existing particle velocities.

**Influence** %: controls how much influence VDB forces will have on particle velocities.

**Variation** %: the per-particle percentage of variation to apply.

#### **Influence**

**X/Y/Z**: controls which axis the forces will affect.

**Relative to mass**: controls whether particle forces will be relative to particle masses.

**Simulate substeps**: forces will be interpolated in a way that simulates the addition of forces at smaller simulation substeps.

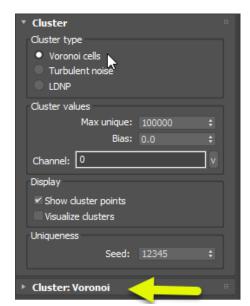
## Uniqueness

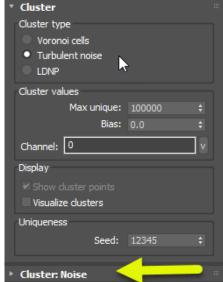
# **Cluster operator**

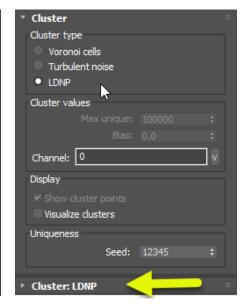


The Cluster operator allows you to partition particles into groups and then save the resulting group index to a custom data channel, for later use in other operators.

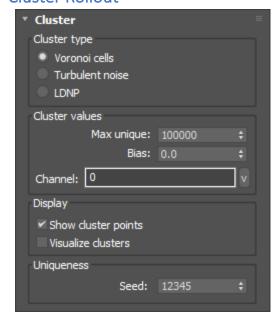
The rollouts change depending on what cluster type is selected and available options also change







## **Cluster Rollout**



#### **Uniqueness**

**Seed**: the seed value for all varied parameters.

## **Cluster Type**

**Voronoi cells**: cluster groups will take the shape of voronoi cells, derived from nearest-point proximity groupings.

**Turbulent noise**: cluster groups will be derived from 3D turbulent noise values.

**LDNP**: cluster groups will be derived from LDNPs (local distributions of neighboring particles). This is an approximated k-means clustering method.

#### **Cluster Values**

**Max unique**: the maximum number of cluster groups.

**Bias**: the amount of bias that group indices will have towards lower values.

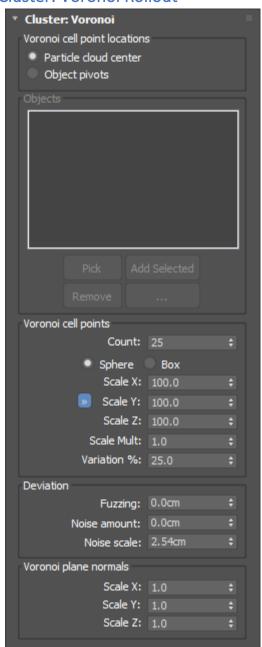
**Channel**: the custom data float channel to save the resulting group indices to.

## **Display**

**Show cluster points**: show the individual cluster points (voronoi cell centers) in the viewport.

**Visual clusters**: display each particle as a random color, based on their cluster group index.

## Cluster: Voronoi Rollout



#### **Voronoi cell point locations**

**Particle cloud center**: the center of the point cloud will be located at the center of the particle cloud.

**Object pivots**: point clouds will be located at the pivot points of input objects.

#### **Objects**

**Input object list**: the list of scene objects to use as point cloud centers.

#### **Voronoi Cell Points**

**Count**: the number of points in the point cloud.

**Sphere/Box**: the shape of the point cloud bounds.

Scale X/Y/Z: the scale of each point cloud axis.

**Scale mult**: the overall scale multiplier for the point cloud.

**Variation %:** the per-particle percentage of variation to apply.

#### **Deviation**

**Fuzzing**: the amount of implicit jitter to add to each particle prior to point-cloud proximity searches, which will blur the borders between the resulting cells.

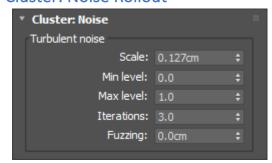
**Noise amount**: the amount of implicit perlin noise to add to each particle prior to point-cloud proximity searches, which will shift the borders between the resulting cells.

**Noise Scale**: the scale of the perlin noise.

## Voronoi plane normals

**Scale X/Y/Z**: the amount of scaling to add to voronoi cell walls, along each axis. Adjusting these values will stretch the overall shape of the resulting voronoi cells.

## Cluster: Noise Rollout



Scale: the scale of the turbulent noise.

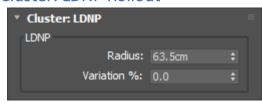
Min level: clamps the minimum value of the resulting noise values.

**Max level**: clamps the maximum level of the resulting noise values.

**Iterations**: the number of iterations of turbulent noise to compute. Higher iterations will result in more detailed noise.

**Fuzzing**: the amount of implicit jitter to add to each particle prior to noise calculations, which will blur the borders between the resulting shapes.

## **Cluster: LDNP Rollout**



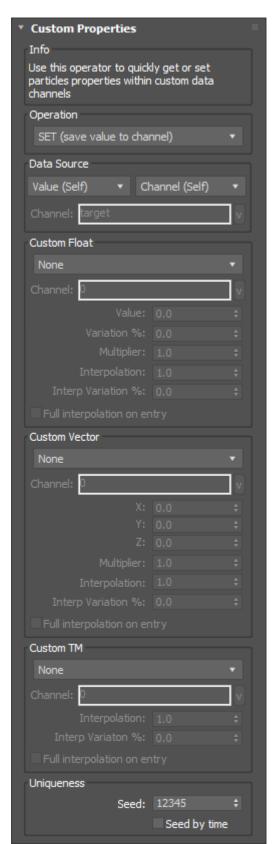
**Radius**: the radius of the particle neighborhoods.

**Variation** %: the per-particle percentage of variation to apply.

# **Custom Properties operator**



The Custom Properties operator provides direct access to each particle's custom data channels.

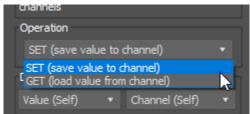


## **Operation**

**Operation list**: the desired operation to perform on custom data.

**SET** (save to channel): Saves the selected particle properties to the selected data channels.

**GET** (**load from channel**): Applies data from the selected channels to the selected particle properties.



#### **Data Source**

**Value Source**: controls whether the values will be the particle's own or the particle's target.

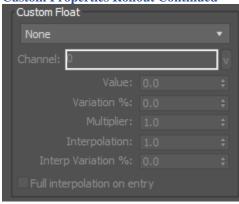
**Channel Source**: controls whether the custom data channel will be the particle's own or the particle's target.

**Target channel**: the name of the custom data channel from which to read the particle's target ID.

### TIP:

The data source options allow you to choose how data flows within the operator. With the introduction of particle targets, it is now possible to read data from a particle and then assign the data to a custom data channel of its target, or vice versa. So, for example, if you wanted a particle to adopt the scale of its target particle, you would have the target particle save its scale to a data channel, and then read that data into the scale value of the source particle. By default, this cross-particle data sharing is disabled within the Custom Properties operator, but can be enabled using these settings.

#### **Custom Properties Rollout Continued**



#### **Custom Float**

Custom float values are individual numerical values composed of single floats.

**Property list**: the particle property to save/load.

**Value**: the absolute float value to save/load.

**Variation** %: the per-particle percentage of variation to apply.

**Channel**: the custom data channel to save/load values to/from.

**Multiplier**: a multiplier applied to the value.

**Interpolation**: the amount to interpolate particle properties from their previous value to the new value.

**Variation %:** the per-particle percentage of variation to apply.

**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

### TIP:

In order to animate a particle property changing into the target value, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.

# **Custom vector**

Custom vector values are Point3 values, composed of x/y/z floats.

**Property list**: the particle property to save/load.

X/Y/Z: the absolute x/y/z values to save/load.

**Channel:** the custom data channel to save/load values to/from.

**Multiplier**: a multiplier applied to the value.

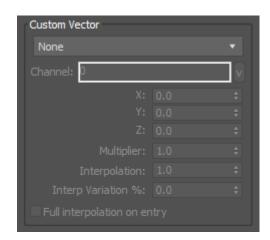
**Interpolation**: the amount to interpolate particle properties from their previous value to the new value.

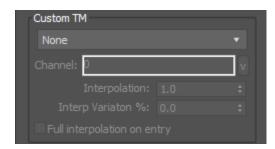
**Variation** %: the per-particle percentage of variation to apply.

**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

# TIP:

In order to animate a particle property changing into the target value, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds





#### **Custom TM**

Custom TM values are Matrix3 values, composed of 4 rows of Point3 values, each composed of x/y/z floats.

**Property list**: the particle property to save/load.

**Channel**: the custom data channel to save/load values to/from.

**Interpolation**: the amount to interpolate particle properties from their previous value to the new value.

**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

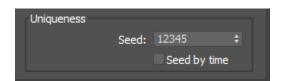
# TIP:

In order to animate a particle property changing into the target value, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.

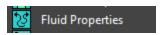
# Uniqueness

**Seed**: the seed value for all varied parameters.

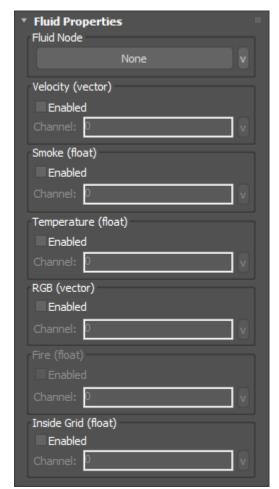
**Seed by time**: the seed value will be incremented with the current time in ticks.



# Fluid Properties operator



The Fluid Properties operator allows you to save various fluid grid properties to particle custom data channels.



#### Fluid Node

**Fluid input object**: the fluid grid object that properties will be loaded from.

## **Velocity (vector)**

**Enabled**: controls whether grid velocity values will be saved to a particle's custom vector data channel.

**Channel**: the custom vector data channel in which to store the data.

## Smoke (float)

**Enabled**: controls whether grid smoke values will be saved to a particle's custom float data channel.

**Channel**: the custom float data channel in which to store the data.

### **Temperature** (float)

**Enabled**: controls whether grid temperature values will be saved to a particle's custom float data channel.

Channel: the custom float data channel in which to store the data.

## Fire (float)

**Enabled**: controls whether grid fire values will be saved to a particle's custom float data channel (FumeFX only).

**Channel**: the custom float data channel in which to store the data.

### Inside (float)

**Enabled**: controls whether a grid inside/outside test will be saved to a particle's custom float data channel.

**Channel**: the custom float data channel in which to store the data.

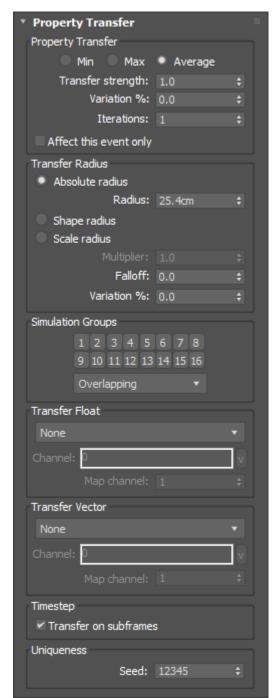
### INFO:

If a particle is inside the fluid grid, a value of '1' will be saved. If a particle is outside the fluid grid, a value of '0' will be saved.

# **Property Transfer operator**



The Property Transfer operator allows particles to spread their properties to neighboring particles.



# **Property Transfer**

**Min**: the minimum value of a property will be transferred among neighboring particles.

**Max**: the maximum value of a property will be transferred among neighboring particles.

**Average**: the average value of a property will be transferred among neighboring particles.

**Transfer strength**: controls the strength of the transfer operation. The lower the value, the less property transfer will occur between neighboring particles.

**Variation %:** the per-particle percentage of variation to apply.

**Iterations**: the number of iterations the transfer operation will run each substep.

**Affect this event only**: properties will only be transferred among particles within the event.

#### **Transfer Radius**

**Absolute radius**: the transfer radius of each particle will be set to a specific value.

**Radius**: the specific transfer radius value.

**Shape radius**: the transfer radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the transfer radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

**Falloff**: the distance beyond the base transfer radius value within which the transfer strength will falloff.

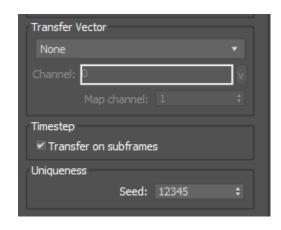
**Variation** %: the per-particle percentage of variation to apply.

#### **Transfer Float**

**Property**: the float property to transfer between neighboring particles.

**Channel**: the custom float channel name to use in custom float transfer mode.

**Map Channel**: the map channel to use in mapping transfer mode.



**Property**: the vector property to transfer between neighboring particles.

**Channel**: the custom vector channel name to use in custom vector transfer mode.

**Map Channel**: the map channel to use in mapping transfer mode.

# **Timestep**

**Transfer on subframes**: when disabled, the operator will only evaluate on whole frames. This can speed up simulations with a small timestep that don't need a lot of transfer fidelity.

# Uniqueness

**Seed**: the seed value for all varied parameters.

# **Script operator**



The Script operator provides advanced control over particles, using C#.

#### INFO:

**tyFlow's** script operator is fast. Combined with its multithreading capabilities, it performs nearly as fast as native c++ code, and is suitable for very complex tasks.

### TIP:

An exhaustive API of all **tyFlow**-related C# functions and parameters can be found within each Script operator's editor window. Users may also reference other C# libraries accessible from Max.NET, like the Math library, for common functions.



## **Script-Accessible Objects**

Script-accessible objects are scene objects that you wish to access directly from within a script.

### INFO:

Scene objects that you wish to access directly from within a script must be registered with the script operator by adding them to the script-accessible object list. Once an object is added to the accessible object list, it can be accessed from within a script using its script-accessible name (displayed in square brackets within the object list). For example, a sphere object added to the list might be displayed as "[obj001] Sphere001", which can then be accessed within the script by the name: obj001.

### **Script-Accessible Texmaps**

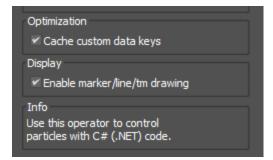
Script-accessible texmaps are scene texmaps that you wish to access directly from within a script.

#### INFO:

Scene texmaps that you wish to access directly from within a script must be registered with the script operator by adding them to the script-accessible texmap list. Once a texmap is added to the accessible texmap list, it can be accessed from within a script using its script-accessible name (displayed in square brackets within the texmap list). For example, a bitmap texture added to the list might be displayed as "[tex001] Map #1", which can then be accessed within the script by the name: tex001.

#### **Optimization**

Cache custom data keys: enables a compile-time optimization which caches custom data channel strings (which are converted into hashmap keys) for faster custom data access.



# INFO:

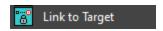
By default, every time a custom data channel is accessed by its channel name, **tyFlow** must convert the string to a hash and then do a hashmap lookup to find the value of that channel for a particular particle. Doing such queries on huge numbers of particles can be very slow. Keeping this optimization enabled will greatly speed up all custom data access, as it manually moves hashmap lookups outside of all particle iteration loops, caching the resulting values for later use.

This should only be disabled for testing purposes.

## **Display**

**Enable marker/line/tm drawing**: enables the drawing of script-based viewport elements (lines, markers, etc). Toggling this setting allows you to choose whether script-based viewport elements will be drawn, without having to modify a script directly (thereby avoiding a simulation reset).

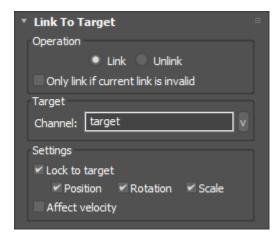
# Link to Target operator



The Link to Target operator allows you to link a particle to its target particle, causing its transform to follow its target particle's transform.

# INFO:

The behavior of the Link to Target operator is similar to regular object linking/parenting within 3ds Max. It is the easiest way to get particles to follow other particles, without introducing heavy solver calculations (like when using Particle or PhysX Bind operators).



# **Operation**

**Link**: particles will be linked to their target.

**Unlink**: any existing links will be broken.

Only link if current link is invalid: linking will only happen if the current link is either invalid (the target particle is deleted or the target ID is invalid) or a link has not been set.

#### **Target**

**Channel**: the channel from which to retrieve the target particle ID.

# **Settings**

**Lock to target**: turning this on will rigidly attach particles to their target. If this setting is off, particles can still move if they have a velocity, and changes to their transform (relative to the target they are linked to) will be made by incrementally measuring target transform changes.

**Position/Rotation/Scale**: controls which portions of a particle's transform will be locked.

**Affect velocity**: when enabled, particle velocities will be set to match the changes they undergo as a result of their link target's motion.

# **Move to Target operator**



The Move to Target operator allows you to move a particle to its target particle's position or to a point on its shape mesh surface.

### INFO:

Unlike the Find Target operator, which moves particles over time by modifying their velocities, the Move to Target operator can be used to move particles instantaneously.



### **Target**

**Channel**: the channel containing the target particle ID.

#### **Move location**

**Target particle position**: the target particle's absolute location in space will be the target point.

Closest point on the target shape mesh: the closest point on the target particle's shape mesh will become the target point.

**NOTE:** If location is set to "closest point on the target shape mesh" and the target particle has no shape mesh, the location mode will revert to "target particle position".

**Target shape mesh vertex pos by index**: the target point will be the vertex on the target particle's shape mesh defined by the vertex index stored in the given custom data float channel.

### **Interpolation**

**Value**: the value used to linearly interpolate between a particle's current location and the target location.

**Variation %**: the per-particle percentage of variation to apply.

#### Offset

**Offset**: the normal-aligned offset to apply to the target location, prior to interpolation.

**Variation** %: the per-particle percentage of variation to apply.

# **Sample**

**Sample**: controls which sampler will be used for closest point determinations.

### **Vertex Index**

**Channel**: the custom float data channel to retrieve the target vertex index value from.

### Uniqueness

**Seed**: the seed value for all varied parameter

# **Set Target operator**



The Set Target operator automates the assignment of the birth ID of one particle to a custom data channel of another.

## **INFO:**

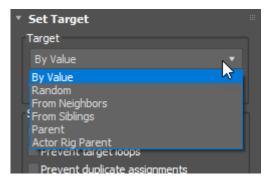
Think of setting a target as a way to tell one particle to focus on another. The "target" is merely the birth ID of another particle in the flow. Many operators allow you to use that target ID in various ways. For example, you can tell particles to bind to their target in a Particle Bind operator, or chase their target in a Find Target operator, etc.

#### TIP:

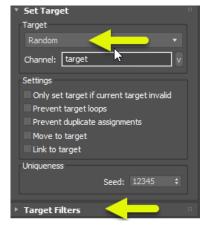
Use a Display Data operator to view which target IDs have been assigned to which particles, or to draw a line between particles and their targets in the viewport.

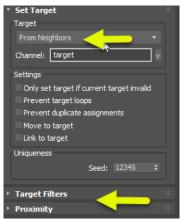
# Set Target Rollout

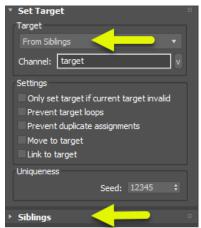
The rollouts available change depending on the target type selected:

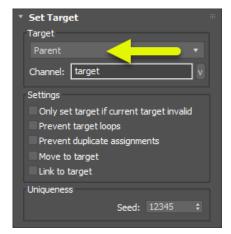




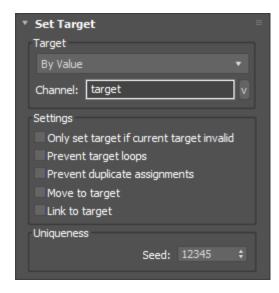












**Mode**: the method to use to assign particle targets.

**Channel**: the custom data float channel to assign target IDs to.

Only set target if current target invalid: particles with a valid (non-deleted) target will not have their target set.

**Prevent target loops**: when enabled, prevents looping particle hierarchies (ie, a particle's target cannot also target it).

**Prevent duplicate assignments**: when enabled, particles will attempt to get a unique target (a target not already assigned to other particles)

**NOTE:** Duplicate prevention happens on a per-step basis. That means if particle A is given a target of 1 at frame 0, then particle B won't get a target of 1 at frame 0. However, if particle A was given a target of 1 at frame 0, that won't prevent particle B from getting a target of 1 at frame 1.

**Move to target**: Moves particles to the location of their target particle.

Link to target: Links particles to their target particle.

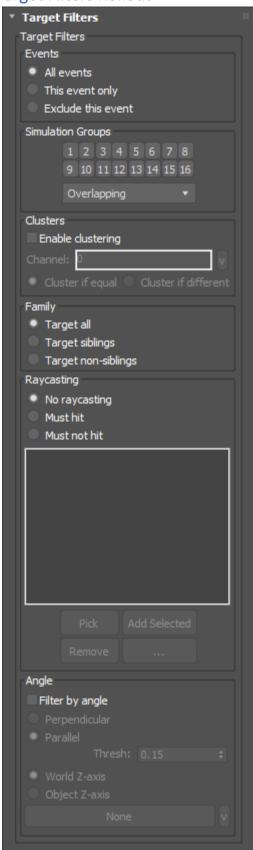
### TIP:

If you need more control over the "move to target" and "link to target" behavior, use the same-named operators instead. These options are merely shortcuts providing fast/simple move and link functionality.

# Uniqueness

**Seed**: the seed value for all varied parameters.

# **Target Filters Rollout**



# Filters Rollout Events

All events: all particles in the flow will be considered for targeting.

**This event only**: only other particles in the event will be considered for targeting.

**Exclude this event**: only other particles in other events will be considered for targeting.

# **Simulation Groups**

**Simulation groups**: controls which particle simulation groups will be targeted.

#### **Clusters**

**Enable clustering**: controls whether target candidates will be determined by cluster values.

**Channel**: the particle data channel to get cluster values from.

Cluster if equal: target candidates must have equal cluster values.

**Cluster if equal**: target candidates must have unequal cluster values.

#### **Family**

**Target all/siblings/non-siblings**: controls whether targets will be filtered by their familial relationships.

# **Raycasting**

**No raycasting/hit/no hit**: controls whether targets will be filtered by a raycast against the defined object list.

**Object list**: the list of objects to raycast against, when raycast filtering is enabled.

# **Angle**

**Filter by angle**: controls whether targets will be filtered by their angle to one another.

**Perpendicular/parallel**: controls whether targets will be filtered based on their perpendicular/parallel angle to a given axis.

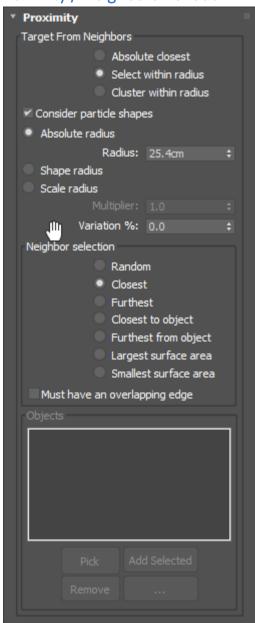
**Thresh**: the threshold to the desired angle within which particles will pass the angle filter.

**World Z-Axis**: the world z-axis (0,0,1) will be the reference vector to which angles will be computed.

**Object Z-Axis**: the local z-axis of a defined object will be the reference vector to which angles will be computed.

**Object**: the object from which to derive the local z-axis.

# Proximity / Neighbors Rollout



**Absolute closest**: the closest particle will become the target candidate for the current particle, regardless of distance.

**Select within radius**: all neighbor particles within a given radius will be target candidates for the current particle.

**Cluster within radius**: all neighbor particles within a given radius will have their target set as the current particle.

**Consider particle shapes**: the vertices of particle shapes will be considered during the neighbor search, allowing for more accurate neighbor detection. If this setting is disabled, only particle positions will be considered.

**Absolute radius**: the neighbor search radius will be an absolute value.

**Radius**: the specific neighbor search radius value.

**Shape radius**: the neighbor search radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the neighbor search radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale neighbor search radius values.

**Variation %:** the per-particle percentage of variation to apply.

# **Neighbor selection**

**Random**: a random particle will be chosen from the list of neighbor candidates as the target.

**Closest**: the closest particle will be chosen from the list of neighbor candidates as the target.

**Furthest**: the furthest particle will be chosen from the list of neighbor candidates as the target.

**Closest to object**: the closest particle to the given scene objects will be chosen from the list of neighbor candidates as the target.

**Furthest from object**: the furthest particle from the given scene objects will be chosen from the list of neighbor candidates as the target.

**Largest surface area**: the particle with the largest surface area will be chosen from the list of neighbor candidates as the target.

**Smallest surface area**: the particle with the smallest surface area will be chosen from the list of neighbor candidates as the target.

#### **Objects**

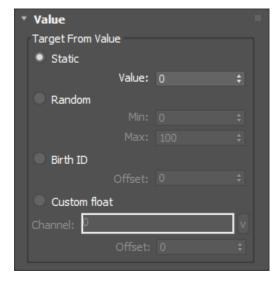
**Object list**: the list of objects used in the relevant closest/furthest modes.

# Siblings Rollout



**Sibling type**: controls which siblings will be considered particle target candidates.

# Value Rollout



**Static**: a specified value will be used as the target ID.

Value: the specified target ID Value.

**Random**: a random value within a min/max range will be used as the target ID.

**Min/Max**: the min/max range from which to choose a random target ID.

**Birth ID**: a permutation of each particle's own birth ID will be used as the target ID.

**Offset**: the amount to offset each particle's own birth ID.

**Custom float**: the target ID will be taken from a custom float data channel.

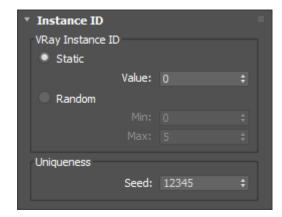
**Channel**: the custom float data channel to take the target ID from.

**Offset**: the amount to offset the target ID value taken from the custom float data channel.

# **Instance ID operator**



The Instance ID can be used to assign a custom Instance ID values to the render instances of applicable particles.



# INFO:

Instance IDs can be utilized by certain VRay texmaps to differentiate between particles independently from material IDs.

### Static

Static mode sets instance IDs to a static value.

Value: the instance ID of particles will be set to this value.

### Random

Random mode chooses a random instance ID from a range of possible IDs.

Min/Max: the range of possible IDs.

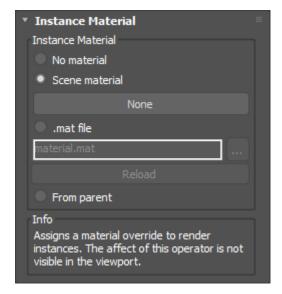
# Uniqueness

**Seed**: the seed value for all varied parameters.

# **Instance Material Operator**



The Instance Material operator can be used to assign a custom material to the render instances of applicable particles.



**NOTE:** The materials applied by this operator **only** affect render instances (for supported renderers like VRay). This operator cannot be used to apply materials in general. In order to assign materials to particles in general, apply the material to the **tyFlow** object itself. If you need multiple materials, assign a Multi/Sub material to the **tyFlow** object itself, and use Material ID operators to differentiate which particles should receive which material.

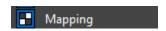
**No material**: clears any material overrides applied to particles.

**Scene Material**: the scene material to apply to render instances of applicable particles.

**.mat file**: the material file whose first material will be applied to render instances of applicable particles.

**From parent**: the material will be inherited from the particle's parent particle.

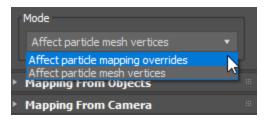
# **Mapping operator**

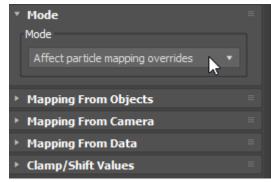


The Mapping operator can be used to assign UVW values to particles.

**NOTE:** The materials applied by this operator **only** affect render instances (for supported renderers like VRay). This operator cannot be used to apply materials in general. In order to assign materials to particles in general, apply the material to the **tyFlow** object itself. If you need multiple materials, assign a Multi/Sub material to the **tyFlow** object itself, and use Material ID operators to differentiate which particles should receive which material.

The available rollouts change depending on the mode selected







# **Mode Rollout**



**Affect particle mapping overrides**: the operator will modify particle mapping override values.

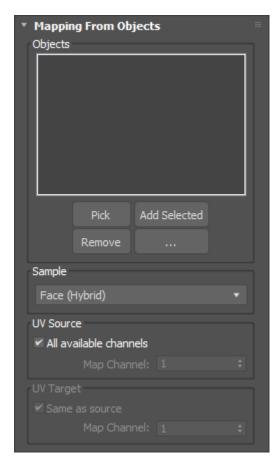
**Affect particle mesh vertices**: the operator will directly modify particle mesh mapping vertex values.

#### INFO:

When in "particle mapping override" mode, each particle will receive a single UVW value for the affected map channels. These override values can be retrieved/changed by other operators, and are not baked directly into the particle meshes until the mesh is sent out from tyFlow (for display, export, rendering, etc). When in "particle mesh vertices" mode, all of the vertices of affected map channels for each particle mesh will receive their own values, which will be baked directly into the particle meshes.

# Mapping From Objects Rollout

By querying closest points on object surfaces, you can assign corresponding mapping values to particles.



**Input object list**: the list of scene objects from which UVW values will be taken.

**Sample type**: controls which sampler will be used to determine closest-object proximities for particles.

#### **UV Source**

**All available channels**: when enabled, UVW values from all available surface channels will be applied to corresponding particle UVW channels.

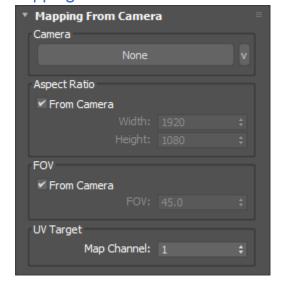
**Map Channel**: the individual map channel to extract UVW values from.

# **UV** Target

**Same as source**: when enabled, the UVW data will be assigned to the same particle UVW channel it was taken from on the source object.

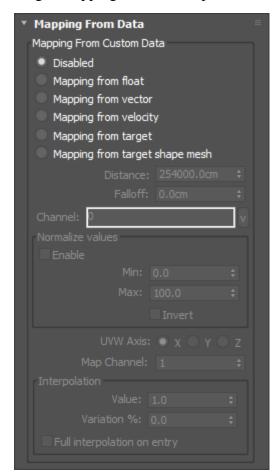
**Map Channel**: the particle UVW channel to assign the UVW data to.

# Mapping From Camera Rollout



# Mapping From Data Rollout

Assigns mapping values from particle custom data channels.



# **Mapping From Custom Data**

**Disabled**: mapping from custom data is disabled.

**Mapping from float**: mapping values will be taken from a particle's custom float data channel.

**Mapping from vector**: mapping values will be taken from a particle's custom vector data channel.

**Mapping from target**: mapping values will be taken from the target particle.

**Mapping from target shape mesh**: mapping values will be taken from the closest point on the target particle's shape mesh.

**Distance/Falloff**: the range to the target particle's shape mesh surface from the particle's location within which UVW values will be retrieved.

**Channel**: the channel to take the custom data value from.

#### Normalize Values

**Enable**: controls whether mapping values will be converted into a value between 0 and 1, based on a target value range.

Min: the minimum value of the range.

**Max**: the maximum value of the range.

**Invert**: inverts the resulting ratio.

**UVW Axis X/Y/Z**: controls which axis of the particle's UVW value to assign custom float data values to.

**Map Channel**: controls which particle mapping channel to assign the UVW value to.

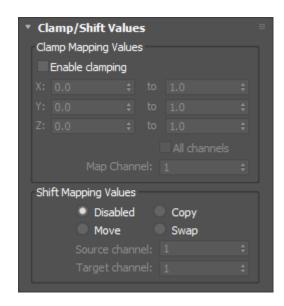
### Interpolation

**Value**: the amount to interpolate particle mapping values from their previous value to the new value.

**Variation** %: the per-particle percentage of variation to apply.

**Full interpolation on entry**: when a particle's event age is zero, its interpolation value will be temporarily set to 1.0.

# Clamp/Shift Values Rollout



# **Clamp Mapping Values**

Allows you to clamp mapping values to specific min/max values.

**Enable clamping**: controls whether clamping will be enabled.

**X/Y/Z**: the min/max x/y/z values to clamp each particle's UVW values between.

**All channels**: clamping will be applied to all mapping channels.

**Map channel**: the specific mapping channel to apply clamping to.

# **Shift Mapping Values**

Allows you to shift particle mapping values between channels.

**Disabled**: no shift operation will be performed.

**Copy**: particle mapping values will be copied from a source map channel to a target map channel. The source channel value will be retained.

**Move**: particle mapping values will be moved from a source map channel to a target map channel. The source channel value will be removed.

**Swap**: particle mapping values will be swapped between a source map channel and a target map channel.

**Source channel**: the source map channel.

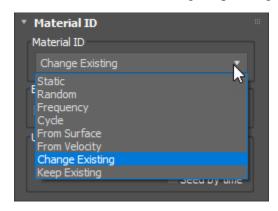
**Target channel**: the target map channel.

# **Material ID operator**

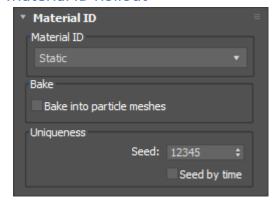


The Material ID operator can be used to assign a material ID to particles.

The available rollouts change depending on the Material ID selection drop down menu



# **Material ID Rollout**



**Operation**: the material ID operation.

**Bake into particle meshes**: instead of being a post-process particle override, the override will be baked into the existing particle mesh and then cleared.

### TIP:

"Keep existing" mode is only useful when "bake into particle meshes" is enabled, since it will leave material ID override values unchanged.

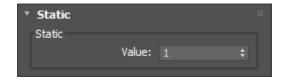
# Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

# Static Rollout

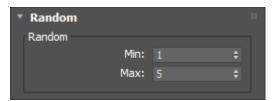
Static mode sets material IDs to a static value.



**Value**: the material ID of particles will be set to this value.

# Random Rollout

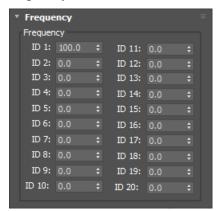
Random mode chooses a random material ID from a range of possible IDs.



Min/Max: the range of possible IDs.

# **Frequency Rollout**

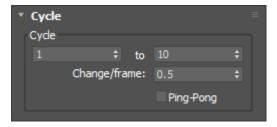
Frequency mode controls the random frequency of specified material IDs.



**ID** #: controls the probability that particles will receive a particular value as their material ID, based on a random sampling of values.

# Cycle Rollout

Cycle mode assigns material IDs to particles each frame, cycling over a range of values as time progresses.



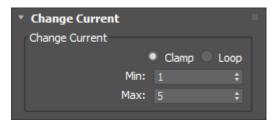
Range: the range of material ID values to cycle over.

**Change/frame**: the amount to progress the cycle each frame.

**Ping-pong**: values will cycle from max to min at the same rate they cycled from min to max.

# **Change Current Rollout**

Change Current mode clamps current material IDs to a range of values.



**Clamp**: clamps the current material ID value to the range using a min/max operation.

**Loop**: loops the current material ID value, to the range using a modulus operator.

Min/Max: the desired range of values

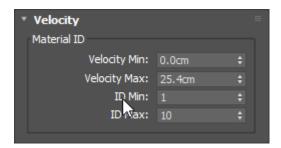
TIP:

In clamp mode, an ID of 7 clamped to a min/max of 1 and 5 will be set to 5: min(7, 5) = 5.

In loop mode, an ID of 7 looped to a min/max of 1 and 5 will be set to 2: 7 % 5 = 2.

# From Velocity Rollout

Velocity mode derives material IDs from particle velocities, by using the ratio between particle velocity magnitudes and a range of target velocity values (which is normalized to a value between 0 and 1). That ratio is then converted to a material ID value between ID min and ID max



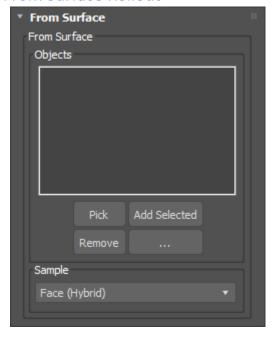
**Velocity Min**: the minimum velocity used to construct the target value range.

**Velocity Min**: the minimum velocity used to construct the target value range.

**ID** Min: the minimum resulting material ID.

**ID** Max: the maximum resulting material ID.

# From Surface Rollout



**Input object list**: the input objects whose surfaces material IDs will be taken from.

**Sample type**: controls which sampler will be used to determine closest-object proximities for particles.

# **Vertex Color operator**

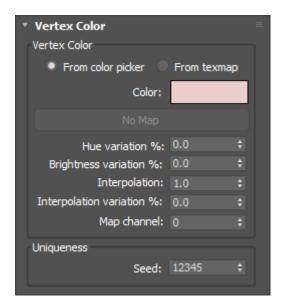


The Vertex Color operator can be used to assign color values to mapping channels of particles.

#### INFO:

This operator is meant for simple color assignments only. If you need more complex behavior (multi-color assignments, color-blending, etc), it's better to use a Custom Properties workflow in combination with a mapping operator, for full control over the values assigned to mapping channels.

**NOTE:** In order to see the color values assigned to particles with this operator in a render, apply a material with a Vertex Color texmap to the **tyFlow** object. To view the vertex color values in the viewport, enable vertex color display in the object properties window of the **tyFlow** object, and add a Mesh operator to your particle with "render only" turned off.



#### **Vertex Color**

**From color picker**: the base color value will be taken from the operator's color picker.

**From texmap**: the base color value will be sampled from the specified texmap, using mapping values assigned to the particles.

**NOTE:** By default, particles have no mapping values assigned to them. You can use a Mapping operator to assign values, or assign them with Custom Properties, etc.

**Hue variation %**: controls the amount of variation to apply to the color's hue.

**Brightness variation %**: controls the amount of variation to apply to the color's brightness.

**Interpolation**: the amount to interpolate particle mapping values from their previous value to the new value.

**Variation** %: the per-particle percentage of variation to apply.

**Map Channel**: controls which particle mapping channel to assign the color values to.

# TIP:

3ds Max's official vertex color channel is typically channel 0, but technically any mapping channel can take color values, so it is not necessary to assign the values to channel 0.

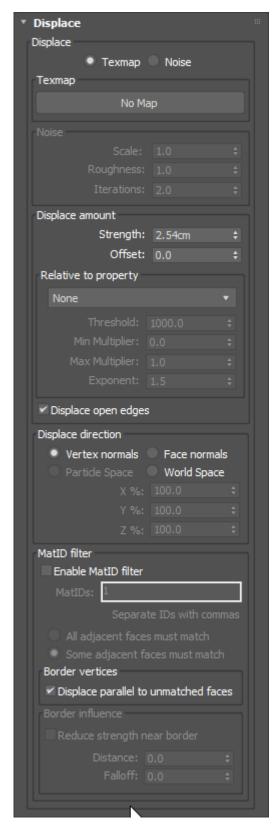
# Uniqueness

**Seed**: the seed value for all varied parameters.

# **Displace operator**



The Displace operator allows you to displace particle mesh vertices using a texture map.



**Texmap/Noise**: controls which displacement method will be used.

#### **Texmap**

In texmap mode, the strength of the displacement is dependent on values derived from a texmap. This mode is dependent on valid mesh UVWs.

**Texmap**: the texture map used to displace particle mesh vertices.

#### Noise

In noise mode, the strength of the displacement is dependent on values derived from a turbulent noise algorithm. This mode is UVW independent.

**Strength**: the strength of the noise.

**Scale**: the scale of the noise.

**Roughness**: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

# **Displace Amount**

**Strength**: the strength of the displacement, in local particle units.

**Offset**: the offset amount to apply to texmap values.

# **Relative to property**

The relative property multiplier will adjust displacement values based on the ratio between the value of a particle property and the specified threshold value.

**Property type**: the particle property to which the threshold value will be compared.



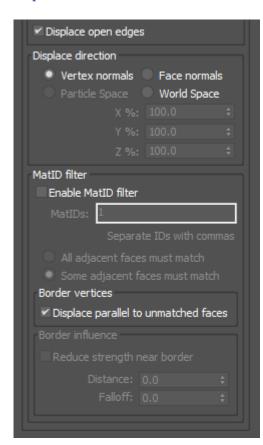
**Threshold**: the target value to which the selected property value of a particle will be compared.

**Min multiplier**: the minimum value of the resulting multiplier.

**Max multiplier**: the maximum value of the resulting multiplier.

**Exponent**: the exponent to which the ratio between the particle property and the selected property will be raised.

### **Displace Rollout Continued**



**Displace open edges**: controls whether vertices on open edges (edges that are not adjacent to two faces) will be displaced.

# **Displace Direction**

**Vertex/Face normals**: vertices will be displaced along the specified mesh normals.

**Particle Space X/Y/Z**: vertices will be displaced along the X/Y/Z planes of the particle's transform.

**World Space X/Y/Z**: vertices will be displaced along the X/Y/Z planes in world-space.

#### MatID filter

MatID filtering allows you to control which vertices in a shape mesh will be displaced.

**Enable MatID Filter**: controls whether material ID filtering will be enabled.

**MatIDs**: the material ID(s) that faces must have in order for their vertices to be displaced.

**All adjacent faces must match**: when enabled, every adjacent face to a vertex must have the matching material ID in order for the vertex to be displaced.

**Some adjacent faces must match**: when enabled, a vertex will be displaced if at least one of its adjacent faces has a matching material ID.

# **Border vertices**

**Displace parallel to unmatched faces**: vertices on the border between a face with a matching material ID and a face without a matching material ID will be extruded in a direction parallel to the planar surface of the face without the matching ID.

#### **Border influence**

**Reduce strength near border**: when enabled, the strength of the displacement will be reduced, depending on how close a displaced vertex is to a border vertex (a vertex on a face whose material ID doesn't match the material ID filter list.

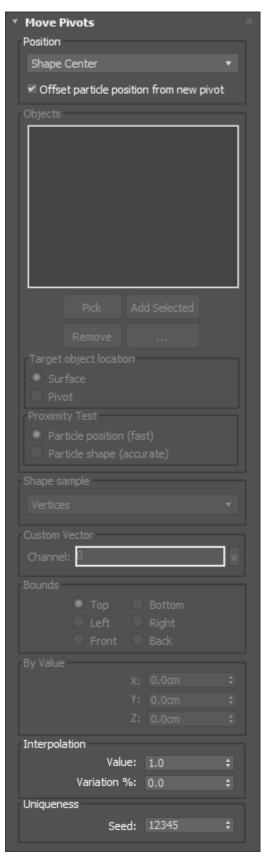
**Distance**: vertices within this distance will not be affected.

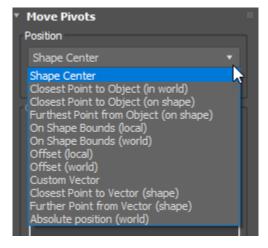
**Falloff**: the effect on vertices beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

# **Move Pivots operator**



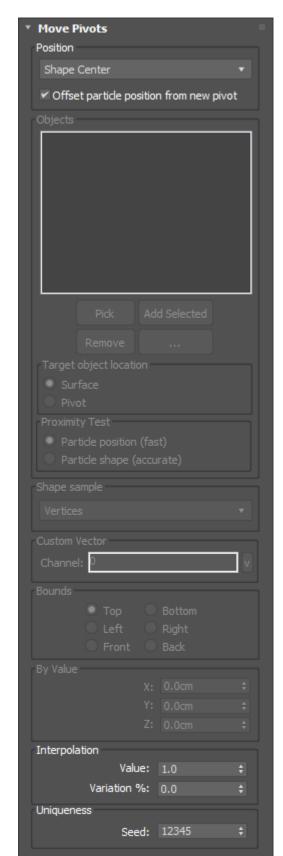
The Move Pivots operator can be used to modify particle pivot points. This is analogous to modifying the pivot of a scene object, if we think of the particle's transform as a pivot and the particle's shape as an object.





#### **Position**

- **Shape center**: moves the particle to the center of its shape mesh.
- Closest point to object (in world): moves the particles to the closest point in world-space to a specified object.
- Closest point to object (on shape): moves the particle to the closest point on its shape to a specified object.
- **Furthest point from object (on shape)**: moves the particle to the furthest point on its shape to a specified object.
- On shape bounds (local): moves the particle to a specified location on its local bounding box.
- On shape bounds (world): moves the particle to a specified location on its axis-aligned bounding box.
- Offset by value (local): moves the particle in the local space of its transform.
- Offset by value (world): moves the particle in world space.
- **Custom Vector**: moves the particle to the location defined by a custom vector.
- Closest point to vector (shape): moves the particles to the closest point to a custom vector on its shape.
- **Furthest point from vector (shape)**: moves the particles to the furthest point from a custom vector on its shape.
- **Absolute position (world)**: moves the particle to an absolute position in world space.



#### **Position**

Offset particle position from new pivot: controls whether the particle position will be offset during the operation, resulting in the relative particle shape location remaining unchanged.

# **Objects**

**Input object list**: the list of input objects.

# **Target object location**

**Surface**: proximity tests will search for the closest point on the input object(s) surface.

**Pivot**: proximity tests will search for the closest input object(s) pivot.

## **Proximity Test**

**Particle Position (fast)**: the closest point to an input object is found by finding the smallest distance between the particle's position and an input object.

**Particle Shape** (accurate): the closest point to an input object is found by finding the smallest distance between the particle's shape and an input object.

# **Shape Sample**

**Vertices**: the vertices of the particle's shape mesh will be used to do the closest/furthest calculations.

**Edge Centers**: the edge centers of the particle's shape mesh will be used to do the closest/furthest calculations.

**Face Centers**: the face centers of the particle's shape mesh will be used to do the closest/furthest calculations.

#### **Custom Vector**

**Channel**: the channel from which to get the custom vector values for the relevant modes.

# Bounds

**Top/Bottom/Left/Right/Front/Back**: the location on the bounding box to move the pivot.

# By Value

**X/Y/Z**: the per-axis values for the applicable modes.

# Interpolation

**Value**: the amount to interpolate particle pivot positions from their previous value to the new value.

**Variation** %: the per-particle percentage of variation to apply.

# Uniqueness

**Seed**: the seed value for all varied parameters.

# **Relax operator**



The Relax operator applies a relaxation algorithm to particle shape meshes.



**Strength**: the strength of the relaxation.

**Iterations**: the number of relaxation iterations to compute.

# **Shape operator**



The Shape operator allows you to assign meshes to particles.

# **Shape Rollout Top Section**



# **Shapes**

**Shape list**: the list of possible shapes that will be assigned to particles.

### TIP:

Select an existing item in the shape list to modify its properties.

**Shape properties** 

### **Distribution**

**Frequency**: the distribution frequency of a particular shape in the list.

### **INFO:**

As particles enter the operator, the shape they are assigned depends on the frequency values of shapes in the list. For example, if one shape has a frequency of 10%, and a second shape has a frequency of 90%, there will be a 10% chance that particles will be assigned the first shape, and a 90% chance that particles will be assigned the second shape.

You do not need to ensure that all shape frequencies add up to 100%. Their values will be normalized automatically if they do not all add up to 100%.

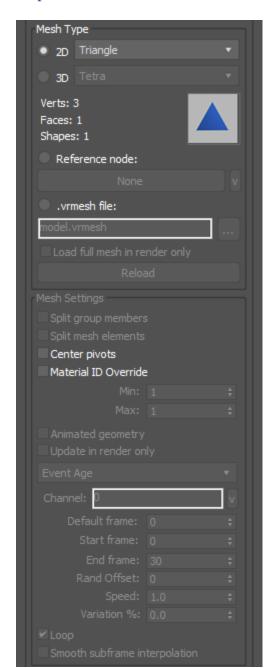
# **Size**

**Override scale**: controls whether to override the scale of particles which are assigned the shape.

**Scale** %: the scale override.

**Variation** %: the per-particle percentage of variation to apply.

#### **Shape Rollout Middle Section**



### **Mesh Type**

**2D**: controls whether the shape will be selected from a list of preset 2D meshes.

**3D**: controls whether the shape will be selected from a list of preset 3D meshes.

**Reference object**: controls whether the shape will be referenced from a scene object in the scene.

**.vrmesh file**: controls whether the shape will be loaded from an existing .vrmesh file.

**Load full mesh in render only**: when enabled, only preview geometry will be displayed for loaded .vrmesh data in the view (this can drastically speed up viewport performance for high resolution proxy files).

**NOTE:** .vrmesh files must be exported with preview geometry enabled, in order to display it in the viewport. Preview geometry settings are available in the standard .vrmesh exporter provided by Chaos Group.

### **Mesh Settings**

**Split group members**: if the shape is referenced from a scene object, and the scene object is a group head, the assigned shape will be chosen from a random child of the group head.

### TIP:

When selecting a group as the reference object, make sure you select the bounding group head object, not a member of the group itself. The group head is denoted in the viewport as a colored bounding box around the group members, and can be displayed by selecting the closed group and choosing Group->Open from the 3ds Max Group menu.

**Split mesh elements**: if the shape mesh has child elements, the assigned shape will be a random child element of the mesh.

**Center pivots**: moves the shape's particle offset to its center.

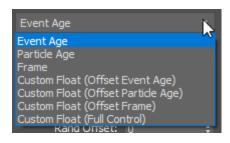
**Material ID Override**: overrides the material ID of the particles the shape is assigned to with a random material ID within a range of specified values.

**Min/Max**: the min/max range of potential material IDs.

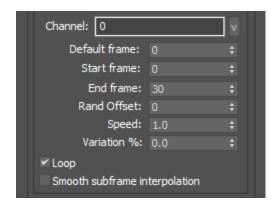
**Animated Geometry**: controls whether animated deformations of the shape will carry over onto particles that are assigned the shape.

**Update in render only**: controls whether animated deformations of the shape are only updated at rendertime. This allows for faster viewport updates.

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## **Middle Section Continued**



# **Animation sync**: controls the offset timing of animated deformations

**Event age**: the start frame of the animation will be synced to the event age of the particle.

**Particle age**: the start frame of the animation will be synced to the age of the particle.

**Frame**: the start frame of the animation will be synced to the current frame of the timeline.

**Custom Float (Offset Event Age)**: the start frame of the animation will be synced to the specified custom float value added to the event age of the particle.

Custom Float (Offset Particle Age): the start frame of the animation will be synced to the specified custom float value added to the age of the particle.

**Custom Float (Offset Frame)**: the start frame of the animation will be synced to the specified custom float value added to the current frame of the timeline.

**Custom Float (Full Control)**: the current frame of the animation will be set to the specified custom float value.

**Default frame**: the default frame that will be loaded in the viewport when "update in render only" is enabled.

**Start frame**: the start frame of the animation to reference.

**End frame**: the end frame of the animation to reference.

**Rand offset**: the per-particle random starting offset within the start/end range of the animation.

**Speed**: the per-particle animation playback speed multiplier.

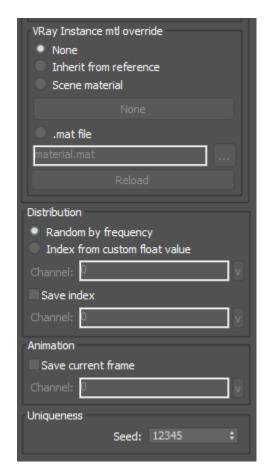
Variation %: the per-particle percentage of variation to apply.

**Loop**: controls whether the animation will loop back to the start frame once the end frame is reached, during sync.

**Smooth subframe interpolation**: controls whether the input geometry will be sampled at subframes or whole frames when playback speed is less than 1.0.

**NOTE:** Enabling smooth subframe interpolation can produce smooth mesh deformations when playback speed is less than 1.0, but can also vastly increase RAM consumption for high resolution geometry when speed variation is greater than 0.0% because the likelihood of two particles being able to share the same mesh is lowered. Enable this setting with caution!

#### **Shape Rollout Bottom Section**



#### **Vray Instance Mtl Override**

**None**: no material override will be assigned to render instances of this shape.

**Inherit from reference**: the material override for render instances of this shape will be taken from the scene object it is referencing.

**Scene Material**: the material override for render instances of this shape will be the specified material.

**.mat file**: the material override for render instances of this shape will be the first material present in the specified material file.

#### **Distribution**

**Random by frequency**: each particle will choose a shape from the list randomly, taking into consideration each list entry's frequency parameter.

**Index from custom float value**: each particle will choose a shape from the list, using a custom data float value as an index to the listbox's entry array.

**Channel**: the custom data float channel to retrieve the index value.

**Save index**: the chosen shape index for each particle will be saved to a custom data float channel.

**Channel**: the custom data float channel to save the index value.

### INFO:

Use "index from custom float value" to specifically choose which shape from the list each particle will be assigned. Doing this will allow you to choose the same shape for each particle across multiple Shape operators in the flow.

### Uniqueness

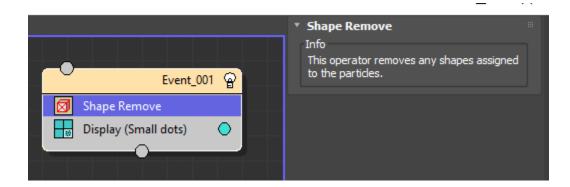
**Seed**: the seed value for all varied parameters.

# **Shape Remove operator**



The Shape Remove operator allows you to remove assigned meshes from particles.

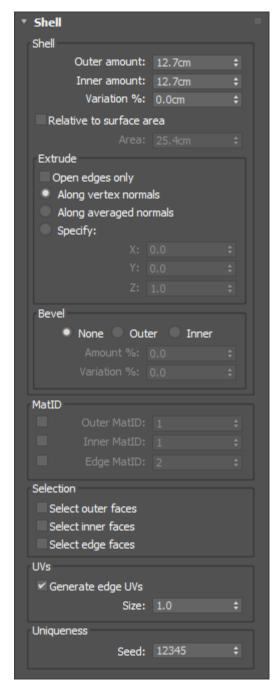
No Parameters or Rollouts for this operator



# **Shell operator**



The Shell operator allows you to give thickness to the faces of particle meshes, similar to how a normal Shell modifier works.



#### Shell

**Outer amount**: the distance to extrude faces along their normal.

**Inner amount**: the distance to extrude faces along the opposite direction of their normal.

**Variation** %: the per-particle percentage of variation to apply.

**Relative to surface area**: the outer/inner amounts will be relative to the overall surface area of the particle divided by the set area.

**Area**: the set area to use in relative mode.

#### Extrude

**Open edges only**: controls whether only elements with open elements are extruded.

**Along vertex normals**: extrudes faces along individual vertex normals.

**Along average normals**: extrudes all faces along the averaged normal direction of the whole mesh.

**Specify X/Y/Z**: extrudes faces in a specified direction.

#### **Bevel**

**None/Outer/Inner**: controls which faces will be beyeled.

**Amount %:** the amount of bevel to apply to the shell faces.

**Variation %**: the per-particle percentage of variation to apply.

#### **MatID**

**Outer MatID Enabled**: controls whether outer faces are given a material ID override.

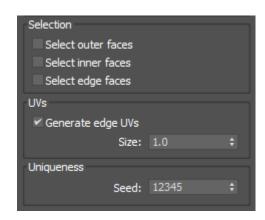
Outer MatID: the override ID to assigned to outer faces.

**Inner MatID Enabled**: controls whether inner faces are given a material ID override.

**Inner MatID**: the override ID to assigned to inner faces.

**Edge MatID Enabled**: controls whether edge faces are given a material ID override.

**Edge MatID**: the override ID to assigned to edge faces.



### Selection

Select outer faces: flags outer faces for selection.Select inner faces: flags inner faces for selection.Select edge faces: flags edge faces for selection.

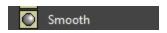
### UVs

**Generate edge UVs**: controls whether UVs will be automatically generated for edge faces.

**Size**: the size of edge UVs.

## Uniqueness

## **Smooth operator**



The Smooth operator assigns smoothing groups to the faces of particle shape meshes.



**Auto Smooth**: When disabled, no mesh faces will share smoothing groups. When enabled, mesh faces will be assigned smoothing groups based on the angles between their adjacent faces.

**Prevent indirect smoothing**: re-processes smoothing groups generated by the auto-smooth function, in order to prevent indirect smoothing.

#### TIP:

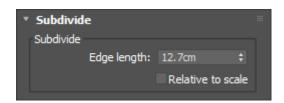
Indirect smoothing happens when adjacent faces (which do not satisfy the angle threshold condition) end up as members of the same smoothing group because they are connected by other adjacent faces which do satisfy the angle threshold condition. "Prevent indirect smoothing" attempts to rectify this problem by shuffling smoothing groups in a way that keeps properly-smoothed adjacent faces as members of the same group, while keeping indirectly-smoothed adjacent faces as members of different groups. Enabling this feature has a performance cost that will cause the auto-smooth algorithm to run slower.

**Threshold**: the angle threshold to use, to compare adjacent faces. If the angle between adjacent faces is less than this value, they will be assigned to the same smoothing group.

## Subdivide operator



The Subdivide operator allows you to subdivide particle meshes.



#### **WARNING:**

Internally, the subdivide operator uses 3ds Max's built-in subdivide modifier for its calculations. This can be slow when applied to particles with hires shape meshes, or when applied to a lot of particles at once. Use with caution.

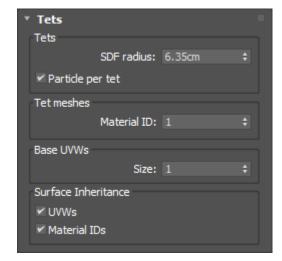
**Edge length**: the target length for all resulting mesh triangle edges.

**Relative to scale**: the edge length value will be relative to each particle's scale.

## **Tets operator**



The Tets operator allows you to convert particle shape meshes into tets.



#### Tets

**SDF radius**: the cell radius of the signed distance field used to create the tet meshes.

**Particle per tet**: when enabled, individual tets will be split out into new, separate particles.

#### Tet meshes

**Material ID**: the material ID to assign to tets, prior to any material ID inheritance.

#### **Base UVWs**

Base UVW coordinates are box-projected coordinates that are assigned prior to any surface inheritance calculations.

**Size**: the real-world size of the box projection.

### **Surface Inheritance**

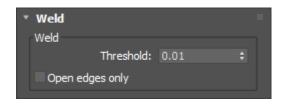
**UVWs**: the outer surface of the generated tets will inherit UVWs from the source mesh.

**Material IDs**: the outer surface of the generated tets will inherit Material IDs from the source mesh.

## **Weld operator**



The Weld operator welds vertices of particle shape meshes.



**Threshold**: the maximum distance threshold between vertices in order for them to be welded together.

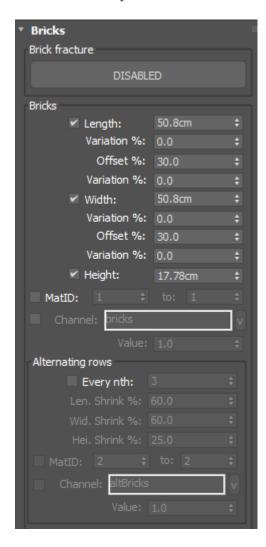
**Open edges only**: only vertices on open edges (edges connected to a single face) will be welded.

## **Brick Fracture operator**



The Brick Fracture operator allows you to slice particle meshes into brick-like patterns.

# Bricks Rollout Bricks Rollout Top Section



## **NOTE:** Each fracture mesh will be turned into a new particle.

#### **Brick Fracture**

**Enable/Disable**: this check button controls whether the brick fracture algorithm will be executed on input particles. It exists as an easy way to enable/disable the algorithm, when tweaking settings (instead of disabling the whole flow to reduce time between setting changes).

#### **Bricks**

**Enable Length**: controls whether slices will be made along the particle's length.

**Length**: the distance between each length-wise slice.

**Variation** %: the per-particle percentage of variation to apply.

**Offset**: the offset applied to length-wise slices, per row.

**Variation** %: the per-particle percentage of variation to apply.

**Enable Width**: controls whether slices will be made along the particle's width.

Width: the distance between each width-wise slice.

**Variation** %: the per-particle percentage of variation to apply.

**Offset**: the offset applied to width-wise slices, per row.

**Variation** %: the per-particle percentage of variation to apply.

**Enable Height**: controls whether slices will be made along the particle's height.

**Height**: the distance between each height-wise slice.

**MatID override**: controls whether a custom material ID will be baked into brick meshes.

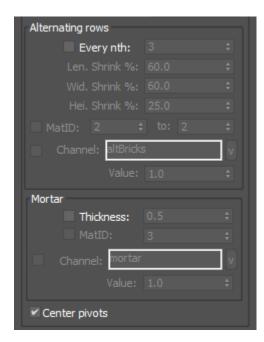
**MatID min/max**: the range of random material ID values to bake into brick meshes.

**Custom Float Data enable**: controls whether a custom float value will be assigned to brick meshes.

**Channel**: the custom float channel where the value will be assigned.

**Value**: the custom float value to assign.

#### **Bricks Rollout Middle Section**



#### **Alternating rows**

Alternating row settings are special shrinkage values you can activate on specific row intervals, to add extra variation to the overall brick pattern.

**Enable alternating rows**: enables alternating row settings.

**Every nth**: controls how often a row will adopt the shrinkage settings.

**Len/Wid/Hei shrink**: controls the scaling factor of length/width/height slices on alternating rows.

**MatID override**: controls whether a custom material ID will be baked into alternating row meshes.

**MatID min/max**: the range of random material ID values to bake into alternating row meshes.

**Custom Float Data enable**: controls whether a custom float value will be assigned to alternating row meshes.

**Channel**: the custom float channel where the value will be assigned.

**Value**: the custom float value to assign.

#### Mortar

When Mortar is enabled, extra slices will be made between bricks, to simulate mortar packed between them.

Mortar enable: enables mortar slices.

Thickness: controls the overall thickness of the mortar effect.

**MatID override**: controls whether a custom material ID will be baked into mortar meshes.

**MatID min/max**: the range of random material ID values to bake into mortar meshes.

**Custom Float Data enable**: controls whether a custom float value will be assigned to mortar meshes.

**Channel**: the custom float channel where the value will be assigned.

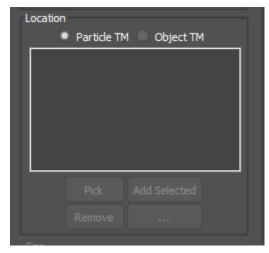
Value: the custom float value to assign.

#### Location

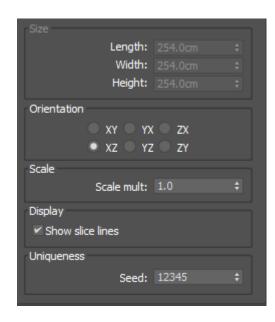
**Particle TM**: the starting point of the slice plane pattern will originate at each individual particle's transform.

**Object TM**: the starting point of the slice plane pattern will originate at specified object transforms.

**Input object list**: the list of input objects to use for slice plane placement.



#### **Bricks Rollout Bottom Section**



#### **Size**

**Length/Width/Height**: the overall size of the slice plane bounding box

#### **Orientation**

**XY/YX/ZX/XZ/YZ/ZY**: the dominant alignment axes of the slice planes.

#### Scale

**Scale mult**: the scale multiplier to apply to the slice planes.

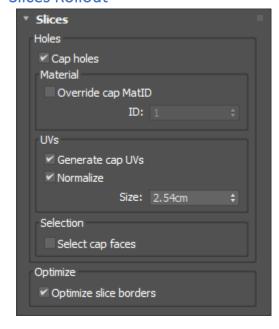
**Show slice planes**: controls whether brick slice planes will be drawn in the view.

**NOTE:** Slice planes will only appear on frames where particles enter the Brick operator for the first time.

#### Uniqueness

• **Seed**: the seed value for all varied parameters.

## Slices Rollout



#### Holes

**Cap holes:** controls whether slices will be capped with new faces.

#### Material

**Override cap MatID**: controls whether cap faces will be given a material ID override.

**ID**: the cap face material ID value.

## UVs

**Generate cap UVs**: controls whether UVW coordinates will be generated on new cap faces.

**Normalize**: controls whether cap UVW coordinates will be normalized.

**Size**: the size of the cap face UVW coordinates.

#### **Selection**

**Select cap faces**: controls whether new cap faces will be flagged for selection.

#### **Optimize**

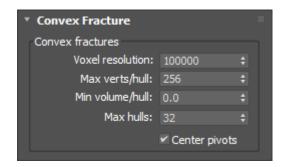
**Optimize slice borders**: controls whether slice borders will be optimized, removing things like extraneous vertices along parallel edges.

## **Convex Fracture operator**



The Convex Fracture operator allows you to fracture particle meshes into groups of convex shapes, using V-HACD.

**NOTE:** Each fracture mesh will be turned into a new particle.



**Voxel resolution**: the resolution setting used to compute the resulting convex hulls.

**Max verts/hull**: controls the maximum number of vertices each resulting hull can have.

**Min volume/hull**: hulls with a volume below this value will be discarded.

**Max hulls**: the maximum number of total hulls returned by the algorithm.

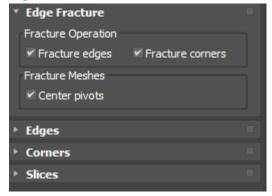
Center pivots: moves each hull's particle offset to its center.

## **Edge Fracture operator**



The Edge Fracture operator allows you to slice particle meshes along the sharp edges of their convex hull.

## **Edge Fracture Rollout**



#### **Fracture operation**

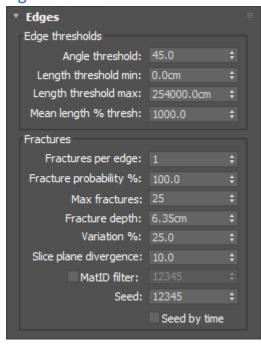
**Fracture edges**: particle meshes will be fractured along the edges of adjacent faces which form a crease.

**Fracture corners**: particle meshes will be fractured at their corners, where groups of edges share a single vertex.

#### **Fracture Meshes**

**Center pivots**: controls whether the pivots of resulting fracture particles will be centered on their mesh.

## **Edges Rollout**



### **Edge Thresholds**

**Edge angle threshold**: the minimum angle between adjacent faces connected to an edge required in order to perform a slice along the edge.

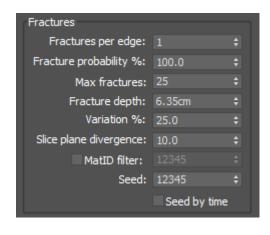
**Length threshold min**: the minimum length an edge must be in order to be fractured.

**Length threshold max**: the maximum length an edge can be in order to be fractured.

Mean length % thresh: an edge's length must not deviate more than this percentage from the average edge length on the mesh, in order to be fractured.

#### TIP:

If the average length of an edge on a mesh is 1.0, and a particular edge length is 1.20 and the mean length % thresh is set to 25%, the edge will be fractured because its length is within a 25% deviation from the average length. However, if the particular edge length is 1.30, it will not be fractured because its length is a 30% deviation from the average length. In other words, the smaller this value, the closer an edge's length must be to the average edge length of the mesh in order to be fractured. Larger values allow for more variation in the lengths of sliceable edges.



**Fractures per edge**: for any given detected edge, the number of initial fractures to generate, prior to any further culling operations.

**Fracture probability**: the probability that any given edge which passed all of the edge threshold tests will be fractured.

**Max fractures**: the maximum number of fractures to compute, per particle.

**Fracture depth**: the distance along the negative edge normal the slice plane will be moved. Higher values result in deeper cuts.

**Variation** %: the per-particle percentage of variation to apply.

**Slice plane divergence**: the maximum angle that any given slice plane normal will diverge from the edge normal.

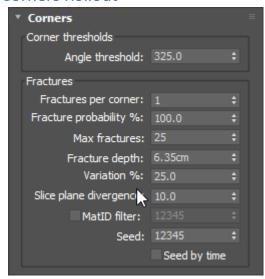
**MatID filter:** when enabled, edges will only be sliced if one or more of their adjacent faces has a material ID matching the specified value.

#### Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

## **Corners Rollout**



**Angle threshold**: the maximum cumulative angle between all edges connected to a single vertex, in order for the vertex to be considered a valid corner.

#### TIP:

The smaller the angle threshold, the sharper a corner must be in order to be fractured.

#### Fractures

**Fractures per corner**: for any given detected corner, the number of initial fractures to generate, prior to any further culling operations.

**Fracture probability**: the probability that any given corner which passed all of the corner threshold tests will be fractured.

**Max fractures**: the maximum number of fractures to compute, per particle.

**Fracture depth**: the distance along the negative corner normal the slice plane will be moved. Higher values result in deeper cuts.

**Variation** %: the per-particle percentage of variation to apply.

**Slice plane divergence**: the maximum angle that any given slice plane normal will diverge from the corner normal.

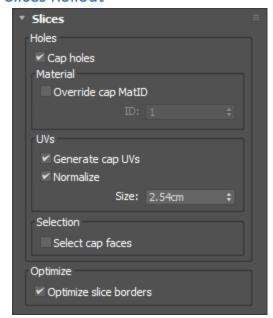
**MatID filter:** when enabled, edges will only be sliced if one or more of their adjacent faces has a material ID matching the specified value.

#### Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks

### Slices Rollout



#### Holes

Cap holes: controls whether slices will be capped with new faces.

#### **Material**

**Override cap MatID**: controls whether cap faces will be given a material ID override.

**ID**: the cap face material ID value.

#### UVs

**Generate cap UVs**: controls whether UVW coordinates will be generated on new cap faces.

**Normalize**: controls whether cap UVW coordinates will be normalized.

**Size**: the size of the cap face UVW coordinates.

#### Selection

**Select cap faces**: controls whether new cap faces will be flagged for selection.

#### **Optimize**

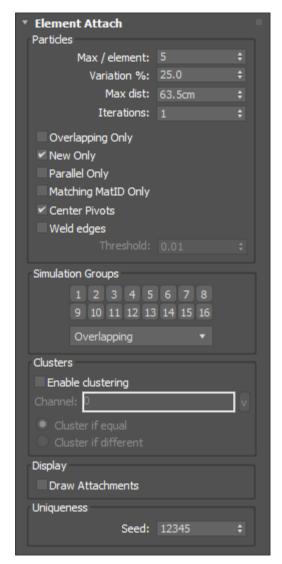
**Optimize slice borders**: controls whether slice borders will be optimized, removing things like extraneous vertices along parallel edges.

## **Element Attach operator**



The Element Attach operator allows you to combine neighboring particle meshes together.

**NOTE:** After a successful attach operation between particles is completed, only the resulting particle (with the newly combined shape mesh) will remain – the rest will be deleted.



#### **Particles**

**Max/element**: controls the maximum number of neighbor particles that can be attached to a particle.

**Variation %:** the per-particle percentage of variation to apply.

**Max dist**: the maximum distance between two particles in order for them to be attach candidates.

**Iterations**: the number of overall attach iterations to perform. This will recursively attach groups of neighboring particles together.

**Overlapping only**: controls whether convex hulls of particles will be computed, in order to determine if particles physically overlap. Only overlapping particles will be attached to each other.

**New only**: when enabled, only particles that are new in the event will be attach candidates.

**Parallel only**: when enabled, only particles that are parallel to each other in space will be attached.

**Matching MatID only**: when enabled, only particles with matching material IDs will be attach candidates for each other.

**Center pivots**: moves each combine mesh's particle offset to its center.

Weld edges: when enabled, the edge vertices of combined meshes will be welded.

**Weld threshold**: the maximum distance between edge vertices of combined meshes, in order for them to be welded together.

#### **Simulation Groups**

**Simulation groups**: controls which particle simulation groups will be candidates for attach.

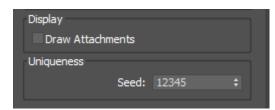
#### Clusters

**Enable clustering**: controls whether attach candidates will be determined by cluster values.

**Channel**: the particle data channel to get cluster values from.

**Cluster if equal**: attach candidates must have equal cluster values.

**Cluster if equal**: attach candidates must have unequal cluster values.



## Display

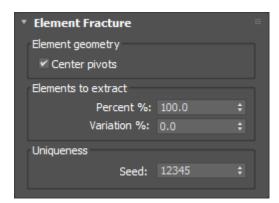
**Draw attachments**: draws a viewport line between successfully attached particles.

## Uniqueness

## **Element Fracture operator**

The Element Fracture operator allows you to fracture particle mesh elements

**NOTE:** Each fracture mesh will be turned into a new particle.



### **Element Geometry**

**Center Pivots**: moves the fractured shape mesh particle offset to its center.

### **Elements to Extract**

**Percent %**: the percentage of shape mesh elements to extract.

**Variation** %: the per-particle percentage of variation to apply.

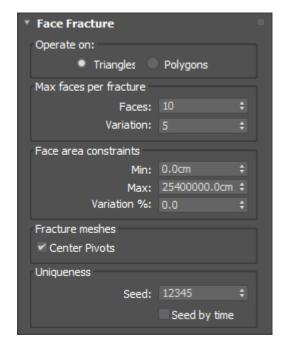
### Uniqueness

## **Face Fracture operator**



The Face Fracture operator allows you to fracture particle mesh faces.

**NOTE:** Each fracture mesh will be turned into a new particle.



#### **Operate On**

**Triangles**: triangles will be treated as single faces.

**Polygons**: polygons (adjacent triangles that share a hidden edge) will be treated as single faces.

#### Max faces per fracture

**Faces**: the number of faces per fracture group.

**Variation**: the per-particle amount of variation to apply.

#### **Face area constraints**

**Min/Max**: the target min/max combined face area per fracture group.

**Variation %**: the per-particle percentage of variation to apply.

#### Fracture meshes

**Center Pivots**: moves the fractured shape mesh particle offset to its center.

#### Uniqueness

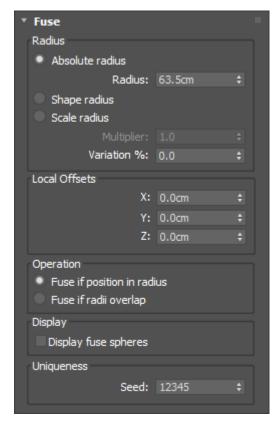
**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

## **Fuse operator**



The Fuse operator uses approximated k-means clustering to fuse neighbor particles within a certain radius together (by deleting all particles except for the particle at the center of each fusion sphere).



#### Radius

**Absolute radius**: the fuse radius of each particle will be set to a specific value.

**Radius**: the specific fuse radius value.

**Shape radius**: the fuse radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the fuse radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

**Variation %**: the per-particle percentage of variation to apply.

#### **Local Offsets**

**X/Y/Z**: position offsets in each particle's local transform space used to control the position of each particle's fusion sphere location.

#### **Operation**

**Fuse if position in radius**: a particle will be fused with another particle if its position is inside the first particle's fuse radius.

**Fuse if radii overlap**: a particle will be fused with another particle if its fuse radius overlaps the first particle's fuse radius.

#### **Display**

**Display fuse spheres**: displays the fuse spheres in the viewport.

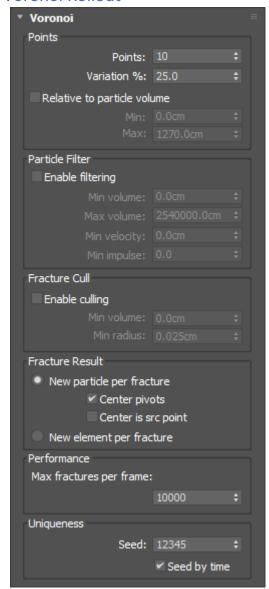
## Uniqueness

## Voronoi Fracture operator



The Voronoi Fracture operator allows you to fracture particle meshes into convex chunks.

#### Voronoi Rollout



#### **Points**

**Points**: the number of points used to create fractures. Roughly the number of resulting fracture meshes, depending on the location of the point cloud.

**Variation** %: the per-particle percentage of variation to apply.

**Relative to particle volume**: controls whether the number of points is relative to the volume of the input shape mesh.

**Min/Max**: the min/max volume values used to create the volume-based point multiplier.

#### **Particle Filter**

**Enable filtering**: enables particle filtering.

**Min volume**: particles with a volume smaller than this value will not be fractured.

**Max volume**: particles with a volume larger than this value will not be fractured.

- \*\* Min velocity\*\*: particles with a velocity below this value will not be fractured.
- \*\* Min impulse\*\*: particles with a PhysX collision impulse below this value will not be fractured.

#### **Fracture Cull**

**Enable culling**: enables fracture culling.

**Min volume**: fracture meshes with a volume below this value will be deleted.

**Min radius**: fracture meshes with a radius below this value will be deleted.

#### **Fracture Result**

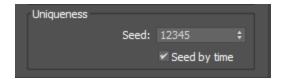
**New particle per fracture**: fracture meshes will be turned into new particles.

**Center Pivots**: moves the fractured shape mesh particle offset to its center.

**New element per fracture**: fracture meshes will be combined as new elements in the particle's shape mesh, replacing the original mesh.

#### **Performance**

**Max fractures per frame**: the maximum number of voronoi fractures to compute, per frame.



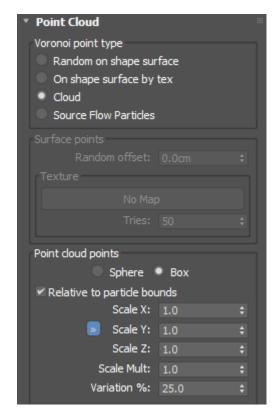
#### Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

## **Point Cloud Rollout**

Top Half of Rollout



#### Voronoi point type

**Random on shape surface**: point cloud points will be generated at random locations on particle mesh surfaces.

On shape surface by tex: point cloud points will be generated on particle mesh surfaces, based on an input texture map.

**Cloud**: point cloud points will be generated at random locations in space.

### **Surface points**

**Random offset**: the amount of random offset to apply to point cloud point positions, along surface normals

**Texmap**: the input texture map to use for particle positions. The sampled intensity will be used as a probability value to determine valid particle locations.

**Tries**: the number of times the texture sampler will try to find a valid particle location before discarding the point cloud point.

#### **Point Cloud Points**

**Sphere/Box**: the bounding shape of the point cloud.

**Relative to particle bounds**: the size of the bounding shape will be relative to the particle's shape size.

Scale X/Y/Z: per-axis scale multipliers for the bounding shape.

**Scale Mult**: a global scale multiplier for the bounding shape.

**Variation** %: the per-particle percentage of variation to apply.

#### Bottom Half of Point Cloud Rollout



#### **Point Cloud Center**

**Particle**: the center of the point cloud will be the particle's position in space.

**Shape center**: the center of the point cloud will be the center of the particle's shape mesh.

**Random in shape bounds**: the center of the point cloud will be a random point within the particle's shape mesh bounding box.

**Random on shape surface**: the center of the point cloud will be a random point on the particle's shape mesh surface.

**Random points on object**: the center of the point cloud will be a random point on an input object's surface.

**At PhysX contact points**: the center of the point cloud will be at a PhysX contact points on the particle's rigidbody hull.

**Input object list**: input objects to use for point cloud positioning.

**Simulation groups**: controls which PhysX collision simulation groups will be candidates for point clouds.

#### Slice plane normals

**Scale X/Y/Z**: the amount of scaling to add to voronoi cell walls, along each axis. Adjusting these values will stretch the overall shape of the resulting voronoi cells.

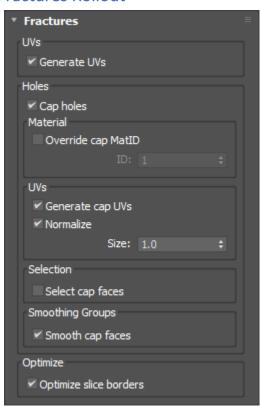
**Show cloud points**: displays point cloud points in the viewport.

#### Uniqueness

**Seed**: the seed value for all varied parameters.

**Seed by time**: the seed value will be incremented with the current time in ticks.

### **Fractures Rollout**



#### **UVs**

**Generate UVs**: controls whether UVW coordinates will be preserved on fractures.

#### Holes

**Cap holes**: controls whether slices will be capped with new faces.

#### **Material**

**Override cap MatID**: controls whether cap faces will be given a material ID override.

**ID**: the cap face material ID value.

#### **UVs**

**Generate cap UVs**: controls whether UVW coordinates will be generated on new cap faces.

**Normalize**: controls whether cap UVW coordinates will be normalized.

**Size**: the size of the cap face UVW coordinates.

#### **Selection**

**Select cap faces**: controls whether new cap faces will be flagged for selection.

#### **Smoothing Groups**

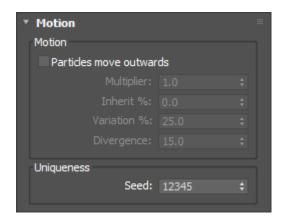
**Smooth cap faces**: controls whether new cap faces will be assigned a smoothing group.

NOTE: When "smooth cap faces" is enabled, the slice algorithm will attempt to assign a smoothing group to cap faces not already assigned to any of the sliced faces around the cap perimeter. If all possible smoothing groups are already used by the sliced faces around the cap perimeter, a smoothing group of 1 will be assigned instead. When "smooth cap faces" is disabled, a smoothing group of 0 will be assigned to cap faces.

#### **Optimize**

**Optimize slice borders**: controls whether slice borders will be optimized, removing things like extraneous vertices along parallel edges.

## **Motion Rollout**



#### **Motion**

**Particles move outwards**: controls whether forces will be applied to fracture particles.

**Multiplier**: the strength of the fracture force to apply to new particles, in the direction of the original particle's position to the new particle's position.

**Inherit** %: the amount of velocity to inherit from the original particle.

Variation %: the per-particle percentage of variation to apply.

**Divergence**: the degrees of divergence to apply to fracture forces.

#### Uniqueness

## **Angle Bind operator**



The Angle Bind operator allows you create binds between adjacent pairs of existing binds which attempt to maintain their angular offset.

**NOTE:** Unlike the Particle Bind operator, this operator does not create bindings between neighboring particles. Instead, it creates binds between the *existing*, *adjacent binds* of a particle. This operator should be applied *after* regular binds are initialized.

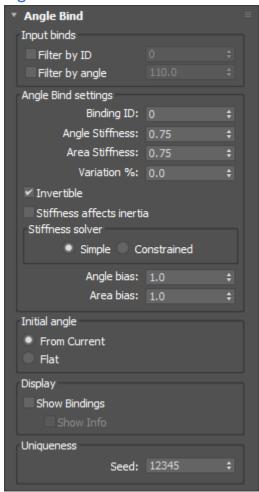
#### TIP:

Existing, adjacent binds are two binds which are connected by a single particle between them. Think of the angle binds generated between them as binds which turn those adjacent binds into a triangle. The adjacent binds are like the top left and right sides of the triangle, and the angle bind is the base of the triangle. The shared particle between the adjacent binds is like the tip of the triangle, and the other two particles at the ends of the adjacent binds are the two points at the base of the triangle. An angle bind attempts to maintain the shape of this implicit triangle created from the existing, adjacent binds

#### INFO:

Bindings are not solved by this operator, only created. Bindings are solved by the global particle bind solver at the end of each simulation step. Bindings created by this operator will persist between events.

## **Angle Bind Rollout**



#### **Input binds**

**Filter by ID**: existing, adjacent binds will be filtered by ID.

**ID**: existing, adjacent binds must have this ID in order to generate an angle bind.

Filter by angle: existing, adjacent binds will be filtered by angle.

**Angle**: two existing, adjacent binds must have at least this large of a relative angle between them, in order to generate an angle bind.

#### **Angle bind settings**

**Binding ID**: the ID to assign to created bindings.

**Angle stiffness**: the stiffness value to assign to the part of the angle bind which attempts to maintain the angle between the existing, adjacent binds.

**Area stiffness**: the stiffness value to assign to the part of the angle bind which attempts to maintain the area of the implicit triangle formed by the existing, adjacent binds.

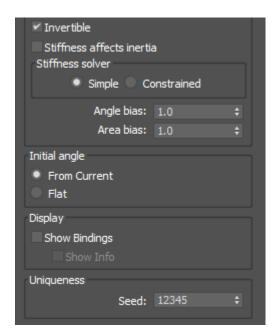
**Variation** %: the per-particle percentage of variation to apply.

**Invertible**: if disabled, the angle bind will attempt to prevent the implicit triangle formed by the existing, adjacent binds from becoming inverted.

#### INFO:

As of tyFlow v0.16083, disabling "invertible" may cause angle binds to become unstable in certain setups. The cause of this instability is still unknown, and will hopefully be improved in future releases.

### **Angle Bind Rollout Continued**



**Stiffness affects inertia**: controls whether the stiffness value of bindings will affect inertia of bound particles.

**NOTE:** "Inertia" is the tendency for an object to resist outside forces. When "stiffness affects inertia" is enabled, binding stiffness will affect how much energy will be transferred between the attached particles. Enabling this setting can help dissipate unwanted energy transfer between particles. This setting has no effect on particles whose bindings have a stiffness value of 1.

#### **Stiffness Solver**

**Simple**: bindings are solved in a straight-forward manner. Their stiffness is mostly dependent on bind solver step count.

**Constrained**: bindings are solved in a way that makes their stiffness somewhat independent of bind solver step count, at a small performance cost. This setting has no effect on bindings whose stiffness value is 1.

**Angle Bias**: the post-initialization multiplier to apply to the rest length of the part of the angle bind which attempts to maintain the angle between existing, adjacent binds.

**Area Bias**: the post-initialization multiplier to apply to the rest length of the part of the angle bind which attempts to maintain the area of the implicit triangle formed by the existing, adjacent binds.

### **Initial Angle**

**From current**: the current angle between the existing, adjacent binds will be used as the target angle of the angle binds.

**Flat**: an angle of 180 degrees will be used as the target angle of the angle binds.

#### TIP:

Choose "flat" when you want the existing, adjacent binds to form a straight line (ex: you are simulating a stiff cord and want to ignore initial angular offsets between existing, adjacent binds).

### **Display**

**Show bindings**: displays bindings as lines in the viewport.

**Show info**: displays binding info (particle ID #1, particle ID #2, stiffness and rest length) in the viewport.

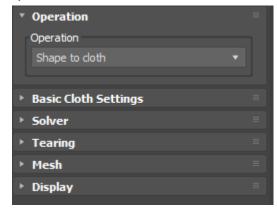
#### Uniqueness

## **Cloth Bind operator**

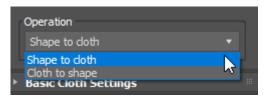


The Cloth Bind operator allows you to convert particle shape meshes into dynamic cloth meshes.

## **Operation Rollout**



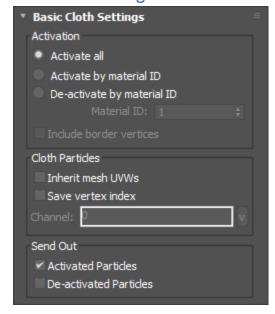
**Operation**: the cloth conversion operation to perform.



**Shape to cloth**: converts a particle shape mesh into a dynamic cloth mesh.

**Cloth to shape**: converts a dynamic cloth mesh into a particle shape mesh.

## **Basic Cloth Settings Rollout**



#### **Send Out**

Activated/de-activated particles: choose which particles will satisfy the operator test condition.

#### **Activation**

Activate all: all new dynamic cloth particles will be activated.

**Activate by material ID**: all new dynamic cloth particles with a matching material ID will be activated.

**De-activate by material ID**: all new dynamic cloth particles will be activated, excluding those with a matching material ID.

**Material ID**: the target material ID to match.

**Include border vertices**: vertices that border faces with differing material IDs will be included as long as one of the material IDs matches the target ID.

#### NOTE:

Activated particles are particles with normal mass. Their position will be affected by the bind solver.

De-activated particles are particles with infinite mass. Their position will not be changed by the bind solver.

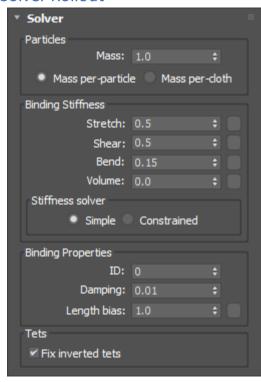
#### **Cloth Particles**

**Inherit mesh UVWs**: causes the new cloth vertex particles to inherit the UVW coordinates of the originating mesh.

**Save vertex index**: saves the cloth mesh vertex index to a custom float data channel in corresponding rig particles.

**Channel**: the custom float data channel to save cloth mesh vertex indices.

## Solver Rollout



#### **Particles**

**Mass**: the mass override for new cloth particles.

**Mass per-particle**: each particle will be assigned the mass override value.

**Mass per-cloth**: each particle will be assigned the mass override value divided by the total vertex count of the cloth mesh.

#### TIP:

The mass per-cloth setting allows you to keep overall cloth masses independent of cloth vertex count, so if you subdivide your input cloth mesh at a later point in time, the overall mass ratio between the cloth as a whole, and other particles, will stay the same

#### **Binding Stiffness**

**Stretch/Shear/Bend/Volume stiffness**: stiffness values for various bindings.

**Stretch/Shear/Bend/Volume texmaps**: the texmaps whose intensity values will be used as a multiplier for the various stiffness values.

#### INFO:

Texmap values are determined by sampling the texmap using the vertices of the input cloth mesh. Since each binding is formed by connecting two vertices, the highest of either texmap value will be used as the multiplier. **INFO:** 

**Stretch bindings**: these bindings are formed between vertices that share a visible edge in the source mesh. They help cloth resist stretching.

**Shear bindings**: these bindings are formed between vertices that share an invisible edge in the source mesh. They also form diagonally between neighbor vertices that span invisible edges. They help cloth resist shearing.

**Bend bindings**: these bindings are formed between any two vertices that are separated by a third vertex between them. They help cloth resist bending.

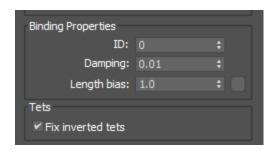
**Volume bindings**: these bindings are formed by casting rays form vertices in the opposite direction of vertex normals, through the interior of a mesh to the other side. They take the shape of internal struts within the mesh and can help cloth maintain overall volume.

#### **Stiffness Solver**

**Simple**: bindings are solved in a straight-forward manner. Their stiffness is mostly dependent on bind solver step count.

**Constrained**: bindings are solved in a way that makes their stiffness somewhat independent of bind solver step count, at a small performance cost. This setting has no effect on bindings whose stiffness value is 1.

#### **Solver Rollout Continued**



#### **Binding Properties**

**ID**: the ID value to assign to bindings.

**Damping**: the amount of post-solve damping to apply to cloth bindings.

**Length bias**: the post-initialization multiplier to apply to binding rest lengths.

**Length bias texmap**: the texmap whose intensity values will be used as a multiplier for the length bias value.

#### **INFO:**

Texmap values are determined by sampling the texmap using the vertices of the input cloth mesh. Since each binding is formed by connecting two vertices, the highest of either texmap value will be used as the multiplier.

#### **INFO**

Since the default bias value is 1.0, the bias texmap multiplier will inversely linearly interpolate values towards 1.0 (for example, an RGB value of [255,255,255] will result in a bias of [length bias spinner value], whereas an RGB value of [0,0,0] will result in bias of [1.0]).

#### **Tets**

**Fix inverted tets**: when simulating cloth composed of tet meshes, the tets can sometimes become inverted (flipping inside out) while undergoing deformations. Enabling this setting will add forces to the simulation at the end of each bind solver which will attempt to prevent/correct tet inversion.

## **Tearing Rollout**



**Enable cloth tearing**: enables cloth tearing.

**NOTE:** Tearing must be enabled in order for **tySlicers** (added within the Particle Break operator) to work on cloth.

#### INFO:

A particle's weakness is its propensity to tear. Weaker particles will cause their attached bindings to tear with less overstretch. Stronger particles require more overstretch before tears will occur. Every time a tear occurs, a progressive weakness multiplier is applied to the attached particles' weakness values. The value of the weakness multiplier can have a large impact on the overall look of the tear. Initial weakness values can be controlled with a texture map, in order to specify where the weakest areas of a cloth are. This gives users more control over tearing behavior.

#### INFO:

Tears caused by tySlicer objects are excluded from tear limits.

#### **Tearing Rollout Continued**



#### Stretch

**Maximum** %: the maximum percentage of stretch a binding may undergo before it will tear, prior to weakness calculations.

**Variation** %: the per-particle percentage of variation to apply.

### Strength

**Tear overstretch frames**: a binding must exceed its maximum stretch percent for these many frames before it will tear.

**Variation** %: the per-particle percentage of variation to apply.

#### Weakness

Use weakness map: controls whether weakness maps are enabled.

**Animated map:** controls whether the weakness map is animated. Animated maps will be updated each time step.

**NOTE:** Depending on the complexity of a cloth simulation, updating weakness values from a texture map can have a measurable performance impact. Therefore, instead of automatically updating a weakness map each frame, users have the ability to select whether or not the map is animated manually. Non-animated maps will only be updated once per simulation.

**Texmap**: the weakness texmap to apply to the cloth, which will affect the weakness of cloth vertices.

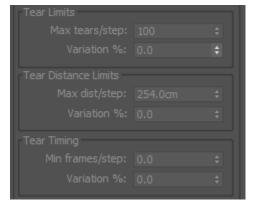
#### **Progressive Weakening**

**Multiplier/tear**: the weakness multiplier to apply to particles that are attached to a torn binding.

**Variation** %: the per-particle percentage of variation to apply.

## **Weakness Limits**

**Min/Max**: the minimum/maximum weakness values a particle may have.



#### **Tear Limits**

**Max tears/strep**: the maximum number of tears that can occur per step.

**Variation** %: the per-particle percentage of variation to apply.

### **Tear Distance Limits**

**Max dist/step**: the maximum cumulative rest length of torn bindings that is allowed per step.

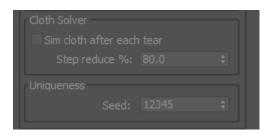
**Variation %**: the per-particle percentage of variation to apply.

#### **Tear Timing**

**Min frames/step**: the minimum duration of frames that must pass before a tear step is processed.

**Variation %:** the per-particle percentage of variation to apply.

## **Tearing Rollout Continued - Bottom Section**



#### **Cloth Solver**

**Sim cloth after each tear**: controls whether additional simulation solves will be performed after each time a bindings is torn.

**Step reduce** %: controls the number of post-tear solver steps will be calculated, as a reduced percentage of the total particle bind solver steps.

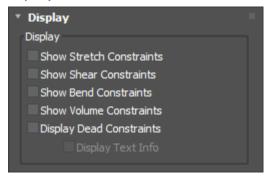
#### TIP:

Every time a binding is torn, it affects the way its neighbor bindings must stretch in order to accommodate the newly torn binding. This change in binding stress can affect which bindings need to be torn next. However, without re-solving all bindings, the tearing algorithm has no way of knowing which areas are experiencing more/less stretching due to the previous tear. By enabling "sim cloth after each tear", the solver has a chance to re-adjust all bindings after each tear which can lead to fewer tearing artifacts. However, re-solving is an expensive operation, and if many tears occur in a single time step this can lead to a significant reduction in solver performance. By increasing the "step reduce %" value, you can decrease the amount of time it takes to perform the intermediate solves while retaining their overall benefit.

#### Uniqueness

**Seed**: the seed value for all varied parameters.

## **Display Rollout**



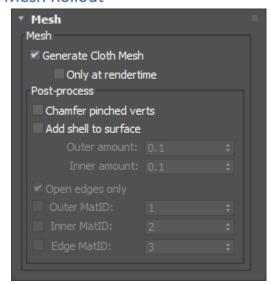
## **Display**

**Show stretch/shear/bend/volume constraints**: controls whether various cloth constraints will be displayed in the viewport.

**Show dead constraints**: controls whether torn constraints will be displayed in the viewport.

**Display text info**: controls whether various pieces of binding debug info will be displayed in the viewport, pertaining to binding particle IDs, rest lengths, stretch ratios, etc.

#### Mesh Rollout



#### Mesh

**Generate Cloth Mesh**: controls whether an actual cloth mesh will be generated by the **tyFlow** object. If disabled, no cloth mesh will be created, even though cloth bindings will still be created between cloth particles.

**At rendertime only**: when enabled, cloth meshes will only be generated during rendertime.

### **Post-process**

Chamfer pinched verts: cloth tears can often lead to degenerate cases where triangles hang from a part of the cloth mesh by a single vertex. Enabling this option will chamfer those single vertex connections between faces, which makes it easier to subdivide the resulting cloth mesh without gaps forming between those faces.

**Add shell to surface**: controls whether the resulting cloth mesh will be given backside/edge faces in order to give it visual thickness.

**Outer amount**: the distance to extrude faces along their normal.

**Inner amount**: the distance to extrude faces along the opposite direction of their normal.

**Open edges only**: controls whether only elements with open elements are extruded.

Outer MatID Enabled: controls whether outer faces are given a material ID override.

**Outer MatID**: the override ID to assigned to outer faces.

**Inner MatID Enabled**: controls whether inner faces are given a material ID override.

**Inner MatID**: the override ID to assigned to inner faces.

Edge MatID Enabled: controls whether edge faces are given a material ID override.

**Edge MatID**: the override ID to assigned to edge faces.

### Cloth Deconversion Rollout



**Delete cloth rig particles**: controls whether cloth rig particles will be deleted, when a dynamic cloth mesh is converted back into a particle shape mesh.

**Carry velocity forward**: controls whether the average velocity of all cloth rig particles will be applied to the resulting cloth shape particle.

## **Cloth Collect operator**



The Cloth Collect operator allows you to collect particles of a cloth mesh that have been spread out into separate events, back into the current event.

There are no parameters for this operator



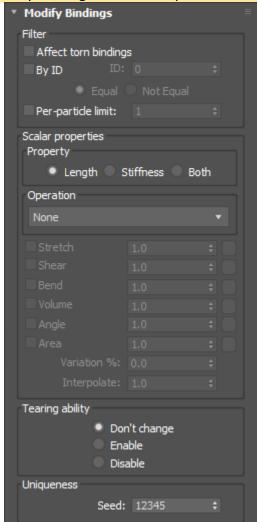
## **Modify Bindings operator**



The Modify Bindings operator allows you to modify the properties of existing particle bindings.

#### TIP:

Only bindings attached to particles influenced by the operator will be affected.



### Filter

**Affect torn bindings**: controls whether bindings that have already been torn will be affected.

**Affect by ID**: controls whether only bindings with a matching ID will be affected.

**ID**: the target binding ID to match.

**Equal**: bindings with an ID that is equal to the target ID will be considered matches.

**Not Equal**: bindings with an ID that is not equal to the target ID will be considered matches.

**Per-particle limit**: when enabled, controls the maximum number of bindings that will be affected, per-particle.

#### **Scalar Properties**

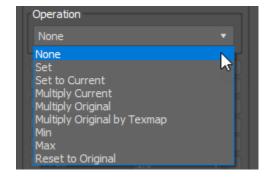
Scalar properties are binding properties defined by a single numerical value.

## **Property**

**Length**: the scalar property being modified will be the binding's rest length.

**Stiffness**: the scalar property being modified will be the binding's stiffness.

**Both**: both length and stiffness will be simultaneously modified.



#### **Operation**

**None**: does not modify scalar properties of bindings.

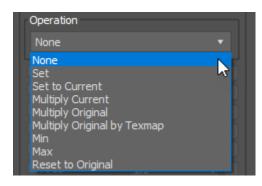
**Set**: changes the scalar property to the specified value.

**Set to current**: changes the scalar property to the current value.

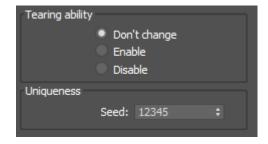
**Multiply current**: multiplies the scalar property by the specified value.

**Multiply original by texmap**: sets the scalar property to its original value, multiplied by the spinner value, interpolated with the texmap mono value.

**Multiply original**: sets the scalar property to its original value, multiplied by the spinner value.







**Min**: sets the scalar property to the specified value, if the scalar property is greater than the specified value.

**Max**: sets the scalar property to the specified value, if the scalar property is less than the specified value.

**Reset to original**: resets the scalar property to its original value at the time the binding was first created.

**Set to current**: changes the rest length value to whatever the current distance between the particles is.

**NOTE:** While cloth bindings are split into different types (stretch/shear/bend/volume), particle bindings are *all* considered stretch bindings.

**Stretch/Shear/Bend/Volume enable**: enables modifications for each type of binding.

**Stretch/Shear/Bend/Volume value**: the specified scalar modifier for each type of binding.

**Stretch/Shear/Bend/Volume texmap**: the texmap used in "multiply by temxap" mode for each type of binding.

**Variation** %: the per-particle percentage of variation to apply.

**Interpolate**: the amount to interpolate particle binding scalar modifications from their previous value to the new value.

#### TIP:

In order to animate particle binding scalar values changing to a particular value over time, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.

**Tearing Ability** 

**NOTE:** Tearing ability must have previously been enabled on bindings at the time of their creation for these settings to take effect.

Don't change: binding tearing ability will not be changed.

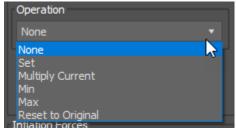
**Enable**: bindings will be flagged as tearable.

**Disable**: bindings will be flagged as un-tearable.

Uniqueness

## Modify Bindings [Cloth] Rollout





#### Tear weakness

These values allow you to directly adjust binding particle tear weakness values.

### **Operation**

**None**: does not modify tear weakness properties of particles.

**Set original**: changes the tear weakness property to the specified value.

**Multiply current**: multiplies the tear weakness property by the specified value.

Min: sets the tear weakness property to the specified value, if the tear weakness property is greater than the specified value.

Max: sets the tear weakness property to the specified value, if the tear weakness property is less than the specified value.

**Reset to original:** resets the tear weakness property to its original value.

#### Values

**Weakness:** The value to use in the tear weakness modification operation.

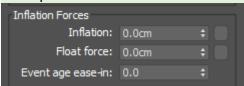
**Variation %:** the per-particle percentage of variation to apply.

**Interpolate**: the amount to interpolate tear weakness modifications from their previous value to the new value.

## TIP:

Operation None

In order to animate binding particle tear weakness values changing to a particular value over time, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.



#### TIP:

Increasing the value of "avent age ease-in" allows you to inflate cloth more gradually

#### **Inflation Forces**

Inflation applies forces along cloth vertex normals during the simulation.

**Inflation**: the strength of the overall inflation force, per particle.

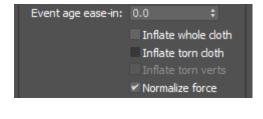
**Inflation texmap**: a texmap that acts as a multiplier on the inflation force value.

**Float**: an artificial force added to the inflation force, in the direction of the world up axis.

Float texmap: a texmap that acts as a multiplier on the float force value.

**Event age ease-in**: the strength of the inflation force will be interpolated from 0 to 1 over this many frames.

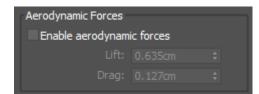
#### **Inflation Forces Continued**











**Inflate torn cloth**: controls whether cloth with torn bindings will inflate.

**Inflate torn verts**: controls whether cloth particles that are directly connected to torn bindings will be inflated.

**Normalize force**: adjusts inflation strength by local face areas.

#### INFO:

If the amount of inflation force applied to each particle controlling a cloth mesh is equal, cloth meshes with asymmetrical topology will inflate in a lop-sided manner. This is because areas of the cloth with a dense number of vertices will receive more overall force than areas with a sparse numbers of vertices. "Normalize force" prevents this from occurring, by adjusting local force values to compensate for topology density.

**Max tear thresh**: If the ratio of torn-to-untorn vertices in a cloth mesh is greater than this value, the cloth will no longer be inflated.

#### TIP:

To simulate air escaping from a torn cloth mesh (which would prevent inflation), adjust the "max tear thresh" to whatever value gives the best result.

**Tear force adjust**: controls the affect tears will have on inflation force strength. The lower this value, the less of an affect tears will have on the amount of inflation force applied.

#### TIP:

The greater the number of overall tears in a cloth mesh, and the greater the "tear force adjust" value, the less of an effect the inflation force will have on vertices. This is an additional way to help simulate air escaping from a torn cloth mesh.

**Display inflation forces**: draws inflation force vectors in the viewport.

#### **Aerodynamic Forces**

Aerodynamic forces are calculated using a thin-foil model of cloth faces. Relative cloth vertex velocities are calculated based on perstep force deltas, and then used to compute lift and drag forces for each cloth vertex.

#### INFO:

Lift and drag force multipliers should generally be kept very small.

**Enable aerodynamic forces**: controls whether lift/drag forces will be calculated.

**Lift**: the lift force multiplier.

**Drag**: the drag force multiplier.

## **Particle Bind operator**

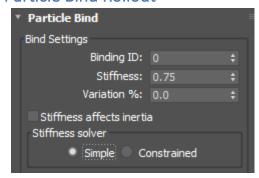


The Particle Bind operator can be used to create bindings between particles.

#### INFO:

Bindings are not solved by this operator, only created. Bindings are solved by the global particle bind solver at the end of each simulation step. Bindings created by this operator will persist between events.

## Particle Bind Rollout



#### **Bind Settings**

**Binding ID**: the ID to assign to created bindings.

**Stiffness**: the stiffness value to assign to created bindings.

**Variation** %: the per-particle percentage of variation to apply.

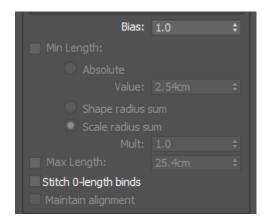
**Stiffness affects inertia**: controls whether the stiffness value of bindings will affect inertia of bound particles.

**INFO:** "Inertia" is the tendency for an object to resist outside forces. When "stiffness affects inertia" is enabled, binding stiffness will affect how much energy will be transferred between the attached particles. Enabling this setting can help dissipate unwanted energy transfer between particles. This setting has no effect on particles whose bindings have a stiffness value of 1.

#### **Stiffness Solver**

**Simple**: bindings are solved in a straight-forward manner. Their stiffness is mostly dependent on bind solver step count.

**Constrained**: bindings are solved in a way that makes their stiffness somewhat independent of bind solver step count, at a small performance cost. This setting has no effect on bindings whose stiffness value is 1.



**Bias**: the post-initialization multiplier to apply to binding rest lengths.

**Enable min length**: applies a minimum length constraint to bindings, during initialization.

**Absolute**: sets the minimum length to an absolute value.

**Absolute value**: the absolute minimum length value.

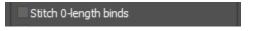
**Shape radius sum**: sets the minimum length to the sum of the attached particles' radii.

**Scale radius sum**: sets the minimum length to the sum of the attached particles' scale values.

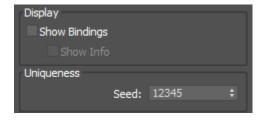
**Mult**: the multiplier applied to shape/scale radius sums.

**Enable max length**: applies a maximum length constraint to bindings, during initialization.

**Max length value**: the absolute maximum length value.







**Stitch 0-length binds**: binds with a rest-length of 0.0 will have their endpoint particles moved to a single position at their center, at the end of the time step.

#### TIP:

Turn on "stitch 0-length binds" if you want to ensure that separate bind hierarchies (ex: cloth) joined together with 0-length binds have no gaps between them. If this option is off, the stretch of the 0-length binds at the end of the time step may be visible.

**Breakable**: controls whether bindings can break, based on various criteria.

%: the percentage of bindings that the break settings will apply to.

**Break by stretch**: enables binding breaking by overstretch.

**Stretch** %: bindings that stretch beyond this percent will break.

Break by compress: enables binding breaking by compression.

**Compress %**: bindings that compress beyond this percent will break.

**Break by min length**: enables binding breaking by explicit minimum length.

\*Min length: bindings that compress to this length will break.

**Break by max length**: enables binding breaking by explicit maximum length.

**Max length**: bindings that stretch to this length will break.

Break by tension: enables binding breaking by tension.

**Tension**: if a binding's attached particles have a tangential velocity magnitude that exceeds this value, the binding will break.

Variation %: the per-particle percentage of variation to apply.

## **Display**

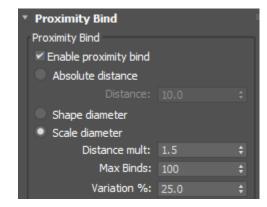
**Show bindings**: displays bindings as lines in the viewport.

**Show info**: displays binding info (particle ID #1, particle ID #2, stiffness and rest length) in the viewport.

#### Uniqueness

## **Proximity Bind Rollout**

Proximity bindings are created between neighbor particles that are a within a threshold distance from each other.



**Enable proximity bind**: controls whether proximity bindings will be created.

**Absolute distance**: controls whether neighbor particles must be within an absolute distance of each other in order to be bound.

**Distance**: the absolute distance value.

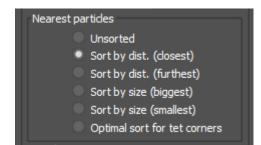
**Shape diameter**: controls whether neighbor particles must be within a distance less than a particle's shape diameter in order to be bound.

**Scale diameter**: controls whether neighbor particles must be within a distance less than a particle's scale diameter in order to be bound.

**Distance mult**: a multiplier applied to the shape/scale diameter.

**Max binds**: the maximum number of bindings that may form between a particle and its neighbors.

**Variation %:** the per-particle percentage of variation to apply.



## **Nearest particles**

When the total number of valid neighbor particles exceeds the "max binds" value, these settings control how to sort the results.

Unsorted: no sorting will occur.

**Sort by dist.** (closest): the neighbors will be sorted from closest to furthest.

**Sort by dist. (furthest)**: the neighbors will be sorted from furthest to closest.

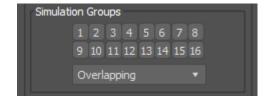
**Sort by size (biggest)**: the neighbors will be sorted from biggest to smallest.

**Sort by size (smallest)**: the neighbors will be sorted from smallest to biggest.

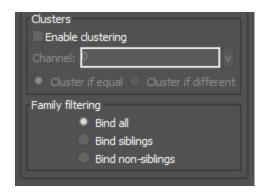
**Optimal sort for tet corners**: Sorts neighbors in an ID-ascending manner, such that all tet corners within range will be bound to each other in a way that is optimal for 0-length bind stitching. When this mode is selected a "max binds" value of 1 is all that is necessary to ensure all tet corner particles will be properly connected.

## **Simulation Groups**

**Simulation groups**: controls which neighbor simulation groups will be bind candidates.



#### Proximity Bind Rollout Continued



#### **Clusters**

**Enable clustering**: controls whether binding candidates will be determined by cluster values.

**Channel**: the particle data channel to get cluster values from.

Cluster if equal: binding candidates must have equal cluster values.

**Cluster if different**: binding candidates must have different cluster values.

## Family filtering

**Bind All**: all particles will be bound, ignoring their relationships.

**Bind siblings**: only sibling particles will be bound to each other.

**Bind non-siblings**: only non-siblings will be bound to each other.

## Family Bind Rollout



**Bind to siblings**: controls whether sibling particles will be bound to each other.

**Bind first and last siblings**: controls whether the siblings at the start and end of a chain of siblings will be bound together, to close the sibling loop.

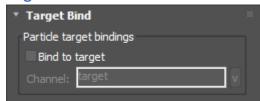
**Proximity**: controls the sibling adjacency level that will be used during the bind operation.

**Bind to parent**: controls whether child particles will be bound to their parent particle.

**NOTE:** Two particles are considered siblings if they were spawned from the same parent particle. Two particles are considered adjacent siblings if they were spawned in order, one immediately before or after the other.

A "proximity" value of 1 means that only adjacent siblings will be bound to each other. A "proximity" value of 2 means that particles will only be bound if they are both adjacent to a third sibling between them. A "proximity" value of 3 means that particles will only be bound if they are both adjacent to two other siblings, who themselves are both adjacent to each other, etc. It essentially relates to the number of sibling adjacency relationships that must exist between two particles, in order for them to be bound.

## **Target Bind Rollout**



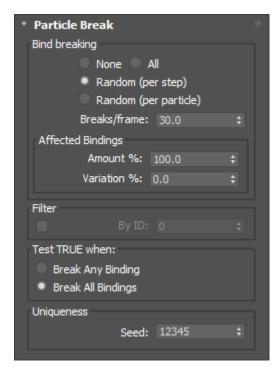
As of 5/5/2020 this rollout is not explained in tyFlow Docs

# Particle Break operator



The Particle Break operator can be used to break bindings between particles.

## Particle Break Rollout



### Uniqueness

**Seed**: the seed value for all varied parameters

# Slicing Rollout



### **Break binding**

**None**: disables bind breaking.

All: breaks all bindings of the event's applicable particles.

**Random** (per step): the "breaks/frame" setting controls how many bindings will be broken per time step, across all applicable event particles.

**Random** (per particle): the "breaks/frame" setting controls how many bindings will be broken per-particle.

**Breaks/frame**: the maximum number of bindings to break per frame.

### **Affected Bindings**

**Amount %:** the percentage of bindings to break.

**Variation** %: the per-particle percentage of variation to apply.

#### **Filter**

**Filter By ID**: controls whether only bindings with a matching ID will be broken.

**ID**: the ID to match.

## **Test TRUE When**

**Break any bindings**: particles will satisfy the test condition when any of their bindings are broken.

**Break all bindings**: particles will only satisfy the test condition when all of their bindings are broken.

The Slicing rollout controls which **tySlicer** objects will be used to slice particle/cloth bindings.

**Input object list**: the **tySlicer** objects to use for bind slicing.

# **Particle Physics operator**

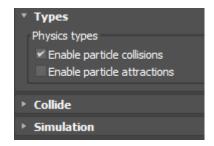


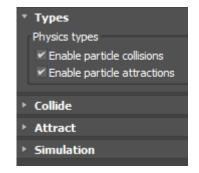
The Particle Physics operator allows you to assign inter-particle collision and cohesion forces.

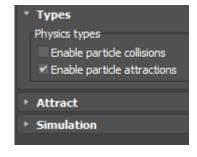
#### INFO:

This operator can be used as the basis for simulations involving sand/grains/snow/etc.

The available rollouts change depending on the physics types selected







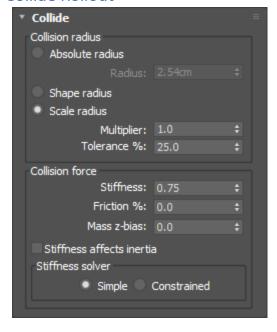
## **Types Rollout**



**Enable particle collisions**: enables inter-particle collision forces.

**Enable particle attractions**: enables inter-particle cohesion forces.

## Collide Rollout



#### **Collision radius**

**Absolute radius**: the collision radius of each particle will be set to a specific value.

Radius: the specific collision radius value.

**Shape radius**: the collision radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the collision radius of each particle will be set to each particle's maximum scale dimension.

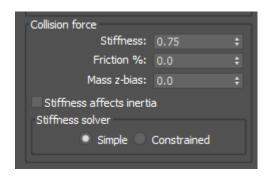
**Multiplier**: the multiplier to apply to shape/scale radius values.

**Tolerance**: a multiplier applied to each particle's radius value, that is used for neighbor searches.

#### NOTE:

The tolerance value controls the size of a buffer around each particle, that allows them to see neighbors which they aren't necessarily touching. A tolerance value greater than zero can help prevent unwanted jittering in the resulting simulation. The greater the tolerance value, the smaller number of time steps you need in a simulation in order to resolve inter-particle collisions. Increasing the tolerance value too much will have a negative impact on performance. A value from 5-50% is ideal.

#### **Collide Rollout Continued**



#### **Collision Force**

**Stiffness**: the strength of the collision resolution force.

**Friction** %: the amount of inter-particle friction that will be applied.

**Mass z-bias**: the amount of artificial mass adjustment to make on stacked particles.

#### NOTE:

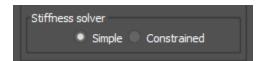
The mass z-bias is a way to artificially decrease the mass of particles stacked on top of each other, in an incremental fashion. This can help stabilize stacks of particles. This value should be kept very low to avoid artifacts - somewhere between .01 and .05 is best.



**Stiffness affects inertia**: controls whether the stiffness value of collisions will affect inertia of colliding particles.

## NOTE:

"Inertia" is the tendency for an object to resist outside forces. When "stiffness affects inertia" is enabled, collision stiffness will affect how much energy will be transferred between the colliding particles. Enabling this setting can help dissipate unwanted energy transfer between particles. This setting has no effect on particles whose collisions have a stiffness value of 1.

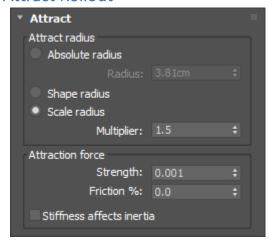


## **Stiffness Solver**

**Simple**: collisions are solved in a straight-forward manner. Their stiffness is mostly dependent on solver step count.

**Constrained**: collisions are solved in a way that makes their stiffness somewhat independent of solver step count, at a small performance cost. This setting has no effect on collisions whose stiffness value is 1.

## **Attract Rollout**



#### **Attract radius**

**Absolute radius**: the attract radius of each particle will be set to a specific value.

**Radius**: the specific attract radius value.

**Shape radius**: the attract radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the attract radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

#### **Attract Force**

**Strength**: the strength of the attraction force.

**Friction %**: the amount of inter-particle friction that will be applied.

**Mass z-bias**: the amount of artificial mass adjustment to make on stacked particles.

**NOTE:** The mass z-bias is a way to artificially decrease the mass of particles stacked on top of each other, in an incremental fashion. This can help stabilize stacks of particles. This value should be kept very low to avoid artifacts - somewhere between .01 and .05 is best.

**Stiffness affects inertia**: controls whether the stiffness value of attraction will affect inertia of attracted particles.

**NOTE:** "Inertia" is the tendency for an object to resist outside forces. When "stiffness affects inertia" is enabled, attraction stiffness will affect how much energy will be transferred between the attracted particles. Enabling this setting can help dissipate unwanted energy transfer between particles. This setting has no effect on particles whose attractions have a stiffness value of 1.

## Simulation Rollout

#### Top Half of Rollout



#### **Integrator**

The integrator is what converts collision/attraction data and assigns it to particle velocities.

**Integrate into bind solver**: controls whether all collision/attraction calculations will be controlled by the Particle Bind Solver, which executes after all operators have been evaluated.

**NOTE:** Unless you need particle physics forces to be evaluated immediately in the operator stack, "integrate into bind solver" should be kept enabled. Disabling "integrate into bind solver" will mean the results of the physics calculations will be immediately available to operators evaluated after the Particle Physics operator.

**OpenCL acceleration**: controls whether the particle neighbor search will be conducted on the GPU.

**Accelerator**: controls which search acceleration structure will be used.

#### INFO:

Hashmap is best when all particles are roughly the same size. KDTree is best when you have a lot of disparity between particle sizes.

**Substeps**: the number of solver steps to perform.

**Affect pos/velocity/both**: controls which particle properties the solver will affect.

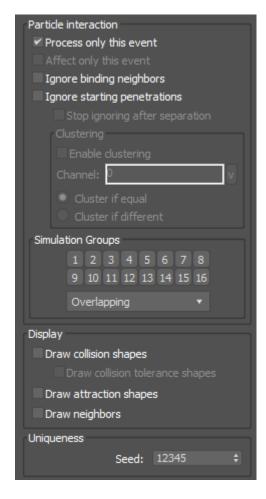
**Affect axis X/Y/Z**: controls which axis of each particle property will be affected.

**Max neighbors**: controls the maximum number of results the neighbor search algorithm will return for a single particle.

**NOTE:** In situations where all particle collision/attraction radii are expected to be non-overlapping and roughly the same size, the "max neighbors" setting can be greatly reduced. The smaller the value, the more memory efficient the GPU algorithm will be. However, if values are too small, or if too many particles are overlapping, or if the difference between the minimum and maximum collision/attraction radii is too large, reducing this value can result in simulation artifacts (jittering, popping, collisions being missed, etc). In those situations, this value may need to be greatly increased.

#### **Simulation Rollout Continued**

#### **Bottom Half of Rollout**



#### **Particle interaction**

**Process only this event**: particle physics forces will only be computed between particles in this event.

**Affect only this event**: particle physics forces will be computed between all relevant particles in the flow, but only particles in the current event will be affected by the forces.

**Ignore binding neighbors**: if two particles are connecting by a binding, particle physics forces will not affect them.

**Ignore starting penetrations**: if two particles are intersecting when they both first enter the event, they will be set to ignore each other.

**Stop ignoring after separation**: if two particles that are ignoring each other (due to starting penetrations) separate, they will be set to no longer ignore each other.

## Clustering

Cluster settings within the "ignore starting penetration" parameters allow you to define which clusters of particles to ignore starting penetrations with.

**Enable clustering**: controls whether ignored starting penetrations will be determined by cluster values.

**Channel**: the particle data channel to get cluster values from.

**Cluster if equal**: particle pairs with starting penetrations must have equal cluster values in order to be ignored.

**Cluster if different**: particle pairs with starting penetrations must have different cluster values in order to be ignored.

## **Simulation Groups**

**Simulation groups**: controls which particle simulation groups will determine which particles will interact.

### **Display**

**Draw collision shapes**: draws spheres representing particle collision radii.

**Draw collision tolerance shapes**: draws spheres representing particle collision tolerance radii.

**Draw attraction shapes**: draws a sphere around each particle, representing its attraction radius.

**Draw neighbors**: draws a connection between all interacting sets of particles.

#### Uniqueness

**Seed**: the seed value for all varied parameters.

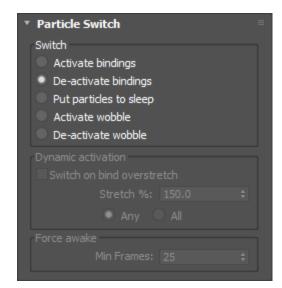
# **Particle Switch operator**



The Particle Switch operator gives you control over which particles will be affected by the particle bind solver.

#### NOTE:

If a particle has bindings attached to it, but you don't want that particle to be affected by the bind solver (for example, you want that particle to stay attached to a static object), you should always be sure to **deactivate** that particle with a Particle Switch operator. When you de-activate a particle, the particle bind solver will treat it as though it has infinite mass, which will allow attached bindings to react to it properly. If you do not de-activate particles properly, your bindings may experience undesirable stretching artifacts.



#### **Switch**

**Activate bindings**: activates particle bindings, causing the bind solver to treat them normally.

**De-activate bindings**: de-activates particle bindings, causing the bind solver to treat them as though they have infinite mass.

**Put particles to sleep:** forces particles to sleep at the current frame.

**Activate wobble**: activates the wobble solver for relevant particles.

**De-activate wobble**: de-activates the wobble solver for affected particles.

## **Dynamic activation**

**Switch on bind overstretch**: causes particles to activate only if attached bindings stretch past a certain point.

**Stretch** %: the threshold stretch percentage that will cause attached particles to activate.

**Any**: activates particles if any of their bindings stretch past the stretch threshold.

**All**: only activates particles if all of their bindings stretch past the stretch threshold.

#### Force awake

**Min Frames**: the minimum number of frames a particle will be forced awake after activation, to prevent premature re-sleeping if the particle's velocity remains under the sleep velocity threshold defined in the Particle Bind Solver settings.

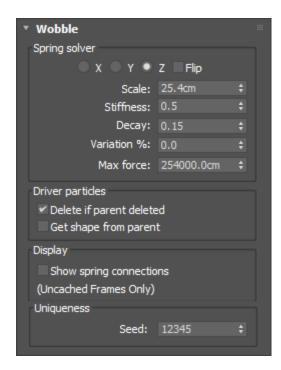
# **Wobble operator**



The Wobble operator allows you to simulate rotational springs on particles.

#### INFO:

The Wobble operator works by coupling each particle with a new child particle in a 2-way relationship that has the following constraints: child particles will attempt to maintain their translational offset from their parent using a simple spring, and parent particles will adjust their orientation to continually look at their child along a specified axis.



### **Spring solver**

**X/Y/Z**: the child offset axis.

**Scale**: the distance each child will be offset from its parent.

**Stiffness**: the stiffness of the spring.

**Decay**: the decay of the spring.

**Variation** %: the per-particle percentage of variation to apply.

**Max force**: the maximum length of the spring force vector.

#### **Driver Particles**

**Delete if parent deleted**: if a parent particle is deleted, the child particle will automatically be deleted.

**Get shape from parent**: controls whether child particles will inherit the shape mesh of their parent.

## **Display**

**Show spring connections**: controls whether a line will be drawn in the viewport connecting child parents with their parent.

#### Uniqueness

**Seed**: the seed value for all varied parameters.

# PhysX Bind operator



The PhysX Bind operator can be used to create bindings between PhysX particles.

#### INFO:

PhysX bindings are not solved by this operator, only created. PhysX bindings are solved by the global PhysX solver at the end of each simulation step. PhysX bindings created by this operator will persist between events.

## PhysX Bind Rollout



## **Type**

**Binding type**: the type of binding to apply to applicable particles

**Distance** (**Tether**): A stiff distance-based binding, that will try to maintain a certain distance between bind points on two PhysX particle rigidbodies.

**Distance (Spring)**: A stretchy distance-based binding, that will try to maintain a certain distance between bind points on two PhysX particle rigidbodies, based on spring/damping settings.

**Rigid** (Glue): a stiff transform-based binding, that will try to maintain perfectly rigid cohesion between PhysX particle rigidbodies.

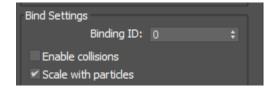
**Rigid (Joint)**: a soft transform-based binding, that will try to maintain some level of cohesion between PhysX particle rigidbodies, based on spring/damping settings.

## TIP:

Sometimes it is better to use Rigid (Joint) bindings with very high spring stiffness, instead of Rigid (Glue) bindings, in order to simulate PhysX particles that are meant to be stuck together. If your Rigid (Glue) bindings enter into a degenerate state that cannot be properly solved, they may introduce more artifacts into a simulation than Rigid (Joint) bindings with very high stiffness.

#### INFO:

Unlike Particle Bind Solver bindings (whose stiffness range is [0-1]), PhysX bindings can take extremely high stiffness values. Stiffness values in the millions are not uncommon for PhysX bindings that need to maintain a high level of stiffness. Don't be afraid to experiment with very large values!

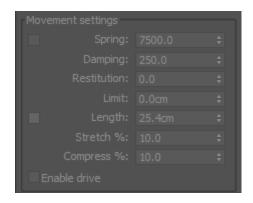


#### **Bind Settings**

**Binding ID**: the ID of created bindings.

**Enable collisions**: controls whether bound PhysX particle rigid bodies should also collide with each other.

**Scale with particles**: controls whether bindings will be automatically adjusted to match the scale of the particles they are attached to, if their scale changes.



## **Movement settings**

**Enable spring**: enables positional springs.

**Spring stiffness**: the stiffness of the springs.

**Damping**: the damping applied to the springs.

**Restitution**: the restitution applied to the springs.

**Limit**: the positional limit of the binding. The solver will attempt to ensure positional deltas do not exceed this value.

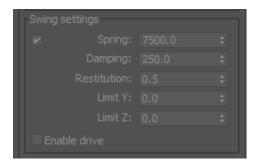
**Length enable**: enables a length-override on the binding.

**Length**: the length-override value.

**Stretch %:** the maximum allowed stretch ratio of the binding.

**Compress** %: the maximum allowed compression ratio of the binding.

**Enable drive**: enables movement drive, which can help to improve the stability of bindings, especially bindings that are very stiff.



## **Swing settings**

Enable spring: enables swing springs.

**Spring stiffness**: the stiffness of the springs.

**Damping**: the damping applied to the springs.

**Restitution**: the restitution applied to the springs.

**Limit Y**: the rotational limit of the binding, around the Y axis.

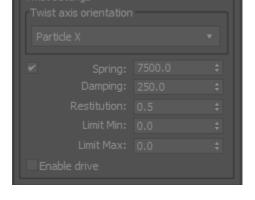
**Limit Z**: the rotational limit of the binding, around the Z axis.

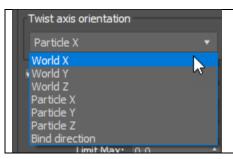
**Enable drive**: enables swing drive, which can help to improve the stability of bindings, especially bindings that are very stiff.



#### Twist axis orientation

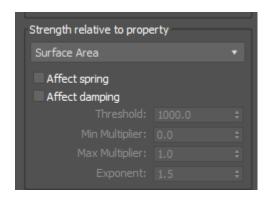
**Axis**: allows you to specify the twist (a-axis) orientation of the bind transform, by choosing which axis to align the twist axis (x-axis) to. The two swing axes will be derived from the specified twist axis orientation.





**NOTE:** Since both swing axes will be derived from the specified twist axis orientation, there is no specific axis orientation setting for them.





**Enable spring**: enables twist springs.

**Spring stiffness**: the stiffness of the springs.

**Damping**: the damping applied to the springs.

**Restitution**: the restitution applied to the springs.

**Limit Min**: the rotational minimum limit of the binding, around the X axis.

**Limit Max**: the rotational maximum limit of the binding, around the X axis.

**Enable drive**: enables twist drive, which can help to improve the stability of bindings, especially bindings that are very stiff.

Strength relative to property

The relative property multiplier will adjust spring/damping values based on the ratio between the value of a particle property and the specified threshold value.

#### TIP:

Imagine a scenario where you're setting up a dynamic tree simulation. You want the binds between the trunk particles to be stronger than the binds between the branch particles (so the trunk is very sturdy, but the branches can sway in the wind). Setting up the values manually on hundreds/thousands of connected particles would be infeasible. Using these settings, you could easily make bind strength relative to particle size, so that larger trunk particles have stronger binds than smaller branch particles.

**Property type**: the particle property to which the threshold value will be compared.

**Affect spring**: controls whether the multiplier will apply to spring values.

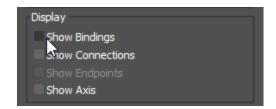
**Affect damping**: controls whether the multiplier will apply to damping values.

**Threshold**: the target value to which the selected property value of a particle will be compared.

**Min multiplier**: the minimum value of the resulting multiplier.

Max multiplier: the maximum value of the resulting multiplier.

**Exponent**: the exponent to which the ratio between the particle property and the selected property will be raised.



#### **Display**

**Show bindings**: bindings between PhysX particle rigidbodies will be drawn in the viewport.

**Show connections**: a connection will be drawn between bound PhysX particles.

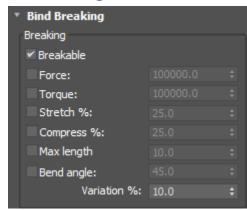
**Show endpoints**: a dot marker will be drawn on either end of the displayed bindings/connections.

### TIP:

If two bound particles are close enough together, their PhysX binding length will be very small - or even zero - so "show bindings" will not be very useful. By choosing "show connections" instead, lines will be drawn between the particles themselves, not simply the locations of their binding's endpoints, so it may be easier to visualize their relationship. Alternatively, you can enable "show endpoints" to display dot markers on either end of display lines, which will help line visibility when line length is close to zero.

**Show Axis**: displays the bind transform axis in the viewport. The twist axis is shown in red, and swing axis are shown in green.

## **Bind Breaking Rollout**



## **Breaking**

**Breakable**: a global override controlling whether any break modes are enabled.

**Force**: a force differential greater than this value between two bound particles will cause their binding to break.

**Torque**: a torque differential greater than this value between two bound particles will cause their binding to break.

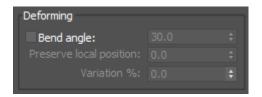
**Stretch %:** bindings that stretch beyond this ratio will break.

**Compress** %: bindings that compress beyond this ratio will break.

**Max length**: bindings that stretch beyond this length will break.

**Bend angle**: bindings that bend beyond this angle will break.

**Variation** %: the per-particle percentage of variation to apply.



#### **Deforming**

**Enable bend deforming**: bindings that bend beyond a certain angle will reset their rest transformation offset.

Bend angle: the bend angle threshold.

**Preserve local offset**: controls the strength of post-deform local bind offset preservation. The lower the value, the more stretching that can occur between bindings as they deform. A value of 1 ensures that only the angle between bindings will deform over time - their relative positions will not be affected.

Variation %: the per-particle percentage of variation to apply.

## **Proximity Bind Rollout**

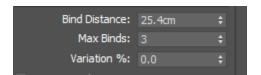


**Particles/Colliders/Ground/World**: controls which rigidbodies of a simulation each particle can be bound to, by proximity.





**Sample type**: controls which sampler will be used to determine closest-surface proximities for particles.



**Bind distance**: the maximum distance between a particle and an enabled binding rigidbody required in order for a new binding to be created between them.

**Max binds**: the maximum number of bindings a particle may form between itself and other relevant rigidbodies.

**Variation** %: the per-particle percentage of variation to apply.

**Improve anchor accuracy**: extra effort will be made to ensure the binding anchor points will be as close together as possible, on the surface of the corresponding rigidbodies.



Improve anchor accuracy

Sort by distance (closest)

Sort by distance (furthest) Sort by volume (biggest)

Sort by volume (smallest)

Sort by surface area (biggest)

Sort by surface area (smallest)

Random

## Rigidbody sorting

When the total number of valid neighbor rigidbodies exceeds the "max binds" value, these settings control how to sort the results.

Unsorted: no sorting will occur.

**Sort by distance (closest)**: the rigidbodies will be sorted from closest to furthest.

**Sort by distance (furthest)**: the rigidbodies will be sorted from furthest to closest.

**Sort by volume (biggest)**: the rigidbodies will be sorted from biggest to smallest by volume.

**Sort by volume (smallest)**: the rigidbodies will be sorted from smallest to biggest by volume.

**Sort by surface area (biggest)**: the rigidbodies will be sorted from biggest to smallest by surface area.

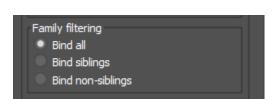
**Sort by surface area (smallest)**: the rigidbodies will be sorted from smallest to biggest by surface area.



## **Simulation Groups**

**Simulation groups**: controls which rigidbody simulation groups will be bind candidates.





#### Clusters

**Enable clustering**: controls whether binding candidates will be determined by cluster values.

Channel: the particle data channel to get cluster values from.

**Cluster if equal**: binding candidates must have equal cluster values.

**Cluster if different**: binding candidates must have different cluster values.

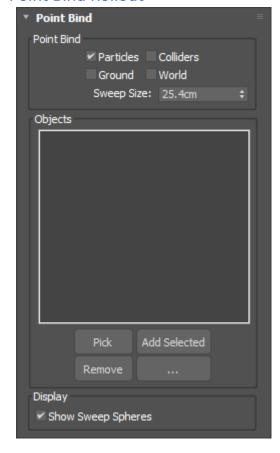
## Family filtering

**Bind All**: all rigidbodies will be bound, ignoring their relationships.

**Bind siblings**: only sibling particles will be bound to each other.

**Bind non-siblings**: only non-siblings will be bound to each other.

## **Point Bind Rollout**



#### **Point Bind**

Point bindings are bindings whose location is controlled by the positions of objects in space.

## TIP:

Proximity bindings will form at the nearest points between two rigidbodies. However, sometimes it is necessary to exert more precise control over the locations of the binding anchor points. Point bindings allow you to control where binding anchor points are set, by positioning a target object in space and ensuring its sweep sphere overlaps the target rigidbodies. Resulting bindings will place their anchor points as close to the target object as possible.

**Particles/Colliders/Ground/World**: controls which rigidbodies of a simulation each particle can be bound to, by points.

**Sweep size**: the sweep size of each target object's sweep sphere.

**NOTE:** Point bindings will only be created between rigidbodies that overlap the target object's sweep spheres.

#### **Objects**

**Input object list**: the list of target objects.

#### **Display**

**Show sweep spheres**: draws the target object sweep spheres in the viewport.

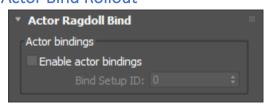
## **Family Bind Rollout**



**Bind to siblings**: controls whether sibling particles will be bound to each other.

**Bind to parent**: controls whether child particles will be bound to their parent particle.

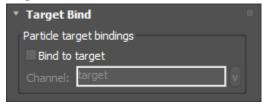
## **Actor Bind Rollout**



**Enable actor bindings**: enables bindings between actor rig particles, whose bind setup ID matches a specified value.

**Bind setup ID**: the bind setup ID to match.

## **Target Bind Rollout**



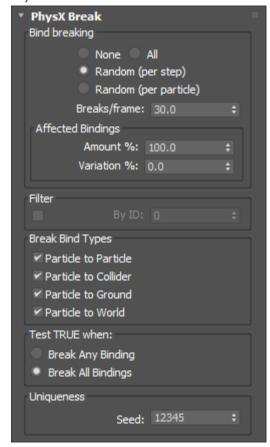
Not Currently Explained in tyFlow Documentation

# PhysX Break operator



The PhysX Break operator can be used to break PhysX bindings.

## PhysX Break Rollout



## **Bind breaking**

**None**: disables PhysX bind breaking.

All: breaks all PhysX bindings of affected particles.

**Random** (per step): the "breaks/frame" setting controls how many PhysX bindings will be broken per time step, across all applicable event particles.

**Random** (per particle): the "breaks/frame" setting controls how many PhysX bindings will be broken per-particle.

**Breaks/frame**: the maximum number of PhysX bindings to break per frame.

#### **Affected Bindings**

**Amount %:** the percentage of PhysX bindings to break.

**Variation** %: the per-particle percentage of variation to apply.

#### **Filter**

**Filter By ID**: controls whether only PhysX bindings with a matching ID will be broken.

**ID**: the ID to match.

#### **Break Bind Types**

[Types]: controls which PhysX binding types will be affected.

#### **Test TRUE When**

**Break any bindings**: particles will satisfy the test condition when any of their PhysX bindings are broken.

**Break all bindings**: particles will only satisfy the test condition when all of their PhysX bindings are broken.

### Uniqueness

**Seed**: the seed value for all varied parameters.

# **Slicing Rollout**



The Slicing rollout controls which **tySlicer** objects will be used to slice PhysX bindings.

**Input object list**: the **tySlicer** objects to use for PhysX bind slicing.

# **PhysX Collision operator**



The PhysX Collision operator allows you to add colliders to a PhysX simulation, and set test conditions for colliding particles.

**NOTE:** Only mesh colliders are supported. Spacewarp colliders are not supported.



#### Colliders

**Input collider objects**: the geometry colliders to add to a PhysX simulation.



## **Hull Type**

**Hull Type**: allows you to choose the collider collision hull.

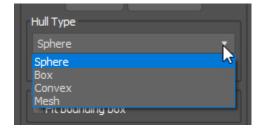
**Sphere**: a perfectly spherical hull.

**Box**: a bounding box hull.

**Convex Hull**: a convex hull encapsulating collider mesh vertices.

Mesh: the collider's mesh will be used as the hull.

**Display Hull**: displays the collider hull in the viewport. **Shape Dimensions** 



Shape Dimensions

Fit bounding box

**Fit Bounding Box**: controls whether spherical hulls will encapsulate the bounding box of the collider's mesh.

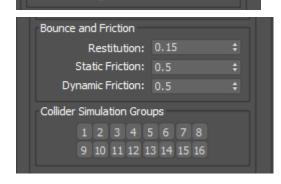
**Bounce and Friction** 

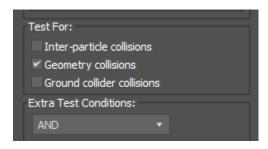
**Restitution**: the restitution coefficient for the collider. Similar to "bounce".

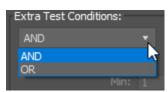
Static/Dynamic friction: friction coefficients for the collider.

**Collider Simulation Groups** 

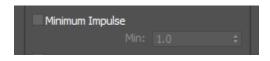
**Simulation groups**: controls which particle simulation groups will be used to filter collisions.

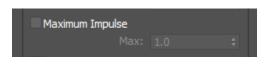




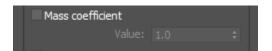












#### **Test For:**

**Inter-particle collisions**: collisions between particles may satisfy the test condition.

**Geometry collisions**: collisions between particles and geometry colliders may satisfy the test condition.

**Ground collider collisions**: collisions between particles and the default PhysX ground plane may satisfy the test condition.

#### **Extra Test Conditions**

If any extra conditions are enabled, they will factor into whether a particle satisfies the test condition.

**Test operator**: the logical operator controlling how extra test conditions are considered.

**AND**: all enabled extra conditions must be satisfied in order for a colliding particle to satisfy the test.

**OR**: any enabled extra conditions must be satisfied in order for a colliding particle to satisfy the test.

## **Collision Count**

A particle must collide a certain number of times before it will satisfy this condition.

Min: the minimum number of collisions.

## **Minimum Impulse**

The minimum impulse of a particle's collision must be above a certain value before it will satisfy this condition.

**Min**: the minimum impulse value.

## **Maximum Impulse**

The maximum impulse of a particle's collision must be below a certain value before it will satisfy this condition.

**Max**: the maximum impulse value.

## **Minimum Penetration**

The penetration value of a particle's collision must be above a certain value before it will satisfy this condition.

**Min**: the minimum penetration value.

#### Mass coefficient

When greater than one, the ratio between the mass of a particle and its collider must be greater than a specified value in order to satisfy this condition. When less than one, the opposite is true. When equal to one, a particle and its collider must have the same mass in order to satisfy this condition.

Value: the test value.

#### Volume coefficient



When greater than one, the ratio between the volume of a particle and its colliding particle must be greater than a specified value in order to satisfy this condition. When less than one, the opposite is true. When equal to one, a particle and its colliding particle must have the same mass in order to satisfy this condition.

Value: the test value.



## **Event Age**

A particle's event age must be within a defined range before it will satisfy this condition.

Min/Max: the event age range.



#### Frame

The current simulation frame must be within a defined range before particles will satisfy this condition.

Min/Max: the frame range.



### **Collider type**

Only particles with a certain PhysX collider rigidbody type will satisfy this condition.

**Dynamic/Kinematic**: the PhysX collider rigidbody type.

## **Simulation Groups**

Only particles matching a certain set of simulation groups will satisfy this condition.

**Simulation groups**: the simulation groups.



# PhysX Fluid operator



The PhysX Fluid operator allows you to simulate two-way interactions between PhysX Shape particle rigidbodies and a PhoenixFD fluid simulation.

## **INFO:**

The PhysX Fluid operator does not rely on PhoenixFD active bodies (available in PhoenixFD v4.x). Instead, it uses a custom PhysX force implementation that is compatible with both PhoenixFD 3.x and PhoenixFD 4.x.

Due to the fact that this operator only uses grid data available in standard PhoenixFD output channels (it does not have access to PhoenixFD's internal FLIP sim data), certain concessions regarding the overall quality of results must be accepted. Influence values will need to be carefully adjusted on a case-by-case basis and may require a decent amount of experimentation to get right.

#### NOTE:

The PhysX Fluid operator calculates forces by sampling interaction points from locations on particle meshes within the specified PhoenixFD fluid grid. When in bounding box mode, the quality of the interaction is affected by how tightly the particle's bounding box conforms to its shape. When in convex hull mode, the quality of the interaction is affected by the number of points sampled on the hull.

#### INFO:

In order to simulate two-way interactions as closely as possible, this operator is designed to run in lock-step with a running PhoenixFD simulation. That means that the flow should not be evaluated before or after the entire PhoenixFD simulation is run, but instead *during* the PhoenixFD simulation process. The easiest way to achieve this is to place the time slider at the starting frame of your PhoenixFD simulation, reset the flow, and then run the PhoenixFD simulation by pressing the "Start" button in the PhoenixFD interface (make sure your **tyFlow** has been added to the list of scene interaction objects within PhoenixFD if the interaction mode is set to "include", and that the particles you wish to interact with PhoenixFD have a proper non-render-only Mesh operator assigned). Alternatively, if you press the "auto-setup" button in the PhysX Fluid operator, the flow will automatically be reset when the PhoenixFD simulation is run (requiring no extra user input in order for the flow to update in lock-step with a running simulation).

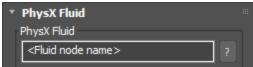
If you need to *restore* a PhoenixFD simulation, move the time slider to the restore frame, reset the simulation manually (letting it re-evaluate up until the restore frame), and then press the "Restore" button in the PhoenixFD interface.

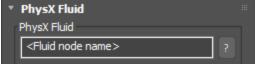
Overall, the basic concept is to simply ensure that the flow updates on a per-frame basis while the PhoenixFD simulation is calculated. Since a running PhoenixFD simulation updates the time slider as it progresses, this will trigger the flow to update at each new frame as well.

## TIP:

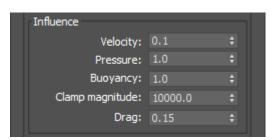
The PhysX Fluid operator applies forces to PhysX particles that take their mass into consideration. Particles with more mass will require higher fluid density or influence values in order to be affected. For better results, ensure particle mass values are proportional to their volumes (this can be done inside a PhysX Shape or Mass operator).

Furthermore, it may be necessary to reduce the **tyFlow's** velocity influence on the PhoenixFD simulation, in order to prevent velocity feedback loops from introducing unwanted motion to the simulation. This can be achieved by reducing the "Motion Velocity Effect" value in the PhoenixFD Properties window for the **tyFlow** object.









**PhoenixFD container name**: The name of the PhoenixFD liquid container object in the scene which will be used to derive fluid forces.

## INFO:

To avoid a circular dependency between tyFlow and PhoenixFD, the PhysX Fluid operator does not hold onto a direct reference to the PhoenixFD node, but instead references it at simulationtype by its name.

Fluid density: the density of the fluid. Higher values will increase the overall influence of the fluid on PhysX particles.

NOTE: By default, the PhysX Shape operator assigns volumerelative mass values to particles. If your particle are large, this could mean they are assigned very large mass values. You can think of the "fluid density" parameter as a parameter which controls the fluid's overall mass. If your particles have a large mass, this value may need to be greatly increased, or you may need to manually assign smaller masses to your PhysX particles in order for the PhoenixFD fluid to influence them.

#### **Influence**

**Velocity**: a multiplier applied to velocity values sampled from the PhoenixFD grid.

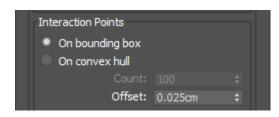
**Pressure**: a multiplier applied to pressure values sampled from the PhoenixFD grid.

NOTE: By default, the PhysX Shape operator assigns volumerelative mass values to particles. If your particle are large, this could mean they are assigned very large mass values. You can think of the "fluid density" parameter as a parameter which controls the fluid's overall mass. If your particles have a large mass, this value may need to be greatly increased, or you may need to manually assign smaller masses to your PhysX particles in order for the PhoenixFD fluid to influence them.

**Buoyancy**: a multiplier applied to a velocity vector pointing straight up, for particle interaction points that are submerged in liquid.

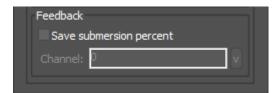
Clamp magnitude: the maximum influence a combined velocity/pressure/buoyancy vector can have on a fully submerged particle. Decrease this value in order to prevent fluid forces from accelerating particles too quickly.

**Drag**: the amount of drag to apply to submerged particles, prior to applying fluid forces.



Display submerged interaction points

Display inactive interaction points



#### **Interaction Points**

On bounding box: fluid interaction points will be generated on the bounding boxes of particle meshes.

On convex hull: fluid interaction points will be uniformly distributed on the convex hulls of particle meshes.

**Count**: the number of interaction points to generate on each particle's convex hull.

**Offset**: the distance from the bounds/hull that the interaction points will be offset.

### TIP:

Adding a slight offset can help ensure interaction points remain submerged in surrounding fluid, when grid cell size is large.

**Display submerged interaction points**: displays interaction points that are submerged in fluid.

**Display inactive interaction points**: displays interaction points that are not submerged in fluid, and are therefore not affected by the grid.

#### **Feedback**

**Save submersion percent**: controls whether the submersion level of a particle (the percentage of its interaction points that are inside of a liquid cell) will be saved to a custom float data channel.

#### TIP:

The submersion percent value ranges from 0-1, depending on how submerged a particle is. 1 means fully submerged.

**Channel**: the custom float data channel where the submersion percent will be saved.

# PhysX Modify operator



The PhysX Modify operator allows you to modify the properties of existing PhysX bindings.

#### INFO:

Only PhysX bindings attached to particles influenced by the operator will be affected. Only **PhysX Joint** and **PhysX Spring** bindings will be affected by stiffness/damping parameters.

## PhysX Modify Rollout



#### **Filter**

**Affect by ID**: controls whether only PhysX bindings with a matching ID will be affected.

**ID**: the target PhysX binding ID to match.

**Equal**: PhysX bindings with an ID that is equal to the target ID will be considered matches.

**Not Equal**: PhysX bindings with an ID that is not equal to the target ID will be considered matches.



## **Scalar Properties**

### **Property**

**Movement**: the properties affecting PhysX binding movement will be affected.

**Swing**: the properties affecting PhysX binding swing will be affected.

**Twist**: the properties affecting PhysX binding twist will be affected.

### **Operation**



**Set**: changes the affected property to the specified value.

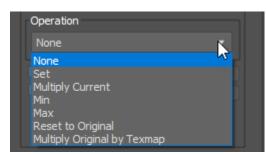
**Multiply current**: multiplies the affected property by the specified value.

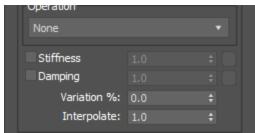
**Min**: sets the affected property to the specified value, if the affected property is greater than the specified value.

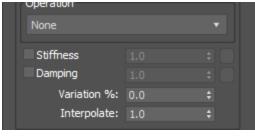
**Max**: sets the affected property to the specified value, if the affected property is less than the specified value.

**Reset to original**: resets the affected property to its original value at the time the PhysX binding was first created.

**Multiply original by texmap**: sets the affected property to its original value, multiplied by the spinner value, interpolated with the texmap mono value.







Stiffness/Damping enable: enables stiffness/damping modification of the affected properties.

Stiffness/Damping value: the specified scalar modifier for each affected property.

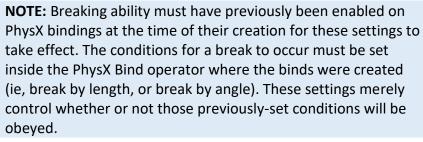
**Variation** %: the per-particle percentage of variation to apply.

**Interpolate**: the amount to interpolate PhysX binding scalar modifications from their previous value to the new value.

## TIP:

In order to animate PhysX binding scalar values changing to a particular value over time, set the operator's timing to "continuous" and set the interpolation value to something less than 1. Smaller interpolation values will result in slower interpolation speeds.

## Breaking



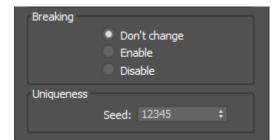
**Don't change**: PhysX bind breaking ability will not be changed.

**Enable**: PhysX bindings will be flagged as breakable.

**Disable**: PhysX bindings will be flagged as un-breakable.

## Uniqueness

**Seed**: the seed value for all varied parameters.

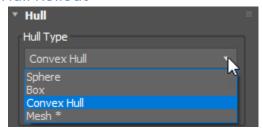


# PhysX Shape operator



The PhysX Shape operator allows you to convert particles into PhysX rigidbodies.

## **Hull Rollout**



### **Hull Type**

**Hull Type**: allows you to choose the particle rigidbody collision bull

**Sphere**: a perfectly spherical hull.

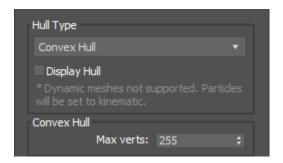
Box: a bounding box hull.

**Convex Hull**: a convex hull encapsulating particle shape mesh

vertices.

**Mesh**: the particle's shape mesh will be used as the hull.

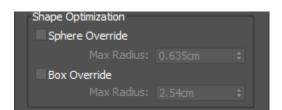
**NOTE:** Dynamic mesh hulls are not supported. Particle rigidbodies with a mesh hull will be set to kinematic.



**Display Hull**: displays the rigidbody hull in the viewport.

**Convex Hull** 

Max verts: the max number of verts allows in the convex hull.



## **Shape Optimization**

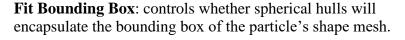
**Sphere Override**: Forces spherical hulls for particles with a radius below a certain threshold.

Max Radius: the radius threshold.

**Box Override**: Forces bounding box hulls for particles with a radius below a certain threshold.

**Max Radius**: the radius threshold.

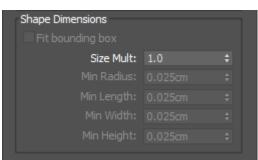




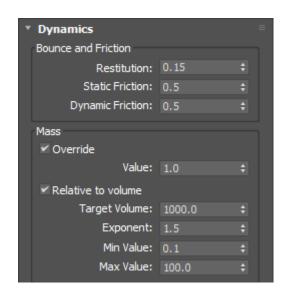
**Size Mult**: the size multiplier for the hull.

Min radius: the minimum radius constraint on spherical hulls.

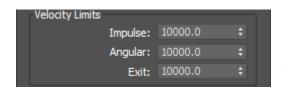
**Min Length/Width/Height**: minimum dimension constraints on box hulls



## **Dynamics Rollout**







#### **Bounce and Friction**

**Restitution**: the restitution coefficient for the rigidbody. Similar to "bounce".

**Static/Dynamic friction**: friction coefficients for the rigidbody.

#### Mass

**Override**: controls whether particle masses will be overridden by this operator.

Value: the desired mass value.

**Relative to volume**: controls whether assigned mass values will be relative to particle shape volume.

**Target volume**: the target volume that particle volumes will be compared to, when assigning mass values.

**Exponent**: the exponent that particle volume ratios will be raised to. Larger exponent values will increase the mass disparity between large and small volumes.

Min Value: the minimum allowed mass value.

Max Value: the maximum allowed mass value.

## **Damping**

**Angular Damping**: the amount of damping to apply to angular (spin) forces acting on rigidbodies.

**Linear Damping**: the amount of damping to apply to linear (movement) forces acting on rigidbodies.

#### TIP:

"Damping" is an effect which reduces the strength of forces over time. Increasing these values will cause particles to spin/move slower over time.

## **Velocity Limits**

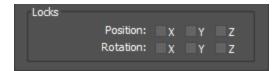
**Impulse**: the maximum impulse (collision) velocity limit for rigidbodies.

**Angular**: the maximum angular (spin) velocity limit for rigidbodies.

**Exit**: the maximum exit (penetration correction) velocity limit for rigidbodies.

## TIP:

Decreasing these values can help prevent jittering and explosive forces that develop as the result of interpenetrations between rigidbodies.

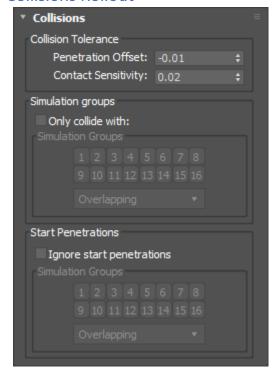


#### Locks

**Position X/Y/Z**: per-axis position locks on rigidbodies.

**Rotation X/Y/Z**: per-axis rotation locks on rigidbodies.

## Collisions Rollout



#### **Collision Tolerance**

**Penetration offset**: the maximum distance that rigidbodies are allowed to penetrate each other.

**Contact sensitivity**: the maximum inflation distance for rigidbody contact detection.

### **Simulation Groups**

Only Collide With: controls whether rigidbodies will only be able to collide with certain simulation groups.

**Simulation groups**: controls which particle simulation groups will be used to filter collisions.

### **Start Penetrations**

**Ignore start penetrations**: controls whether initial penetrations between rigidbodies will be ignored.

**Simulation groups**: controls which particle simulation groups will be used to filter starting penetrations.

## TIP:

If the initial state of your simulation features overlapping rigidbodies, "ignore start penetrations" can prevent explosive forces from being added to the system in order to resolve those penetrations.

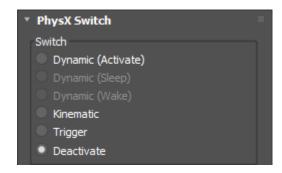
#### INFO:

Penetration resolution between initially-interpenetrating rigidbodies will resume once the rigidbodies fully separate.

# PhysX Switch operator



The PhysX Switch operator gives you control over how PhysX rigidbodies will be treated by the PhysX solver.



#### **Switch**

**Dynamic** (**Activate**): particle PhysX shapes will be treated as dynamic rigidbodies. Affected rigidbodies will be activated.

**Dynamic** (**Sleep**): particle PhysX shapes will be treated as dynamic rigidbodies. Affected rigidbodies will be put to sleep.

**Dynamic** (Wake): particle PhysX shapes will be treated as dynamic rigidbodies. Affected rigidbodies will be awoken.

**Kinematic**: particle PhysX shapes will be treated as kinematic rigidbodies.

**Trigger**: particle PhysX shapes will be treated as trigger rigidbodies.

**Deactivate**: particle PhysX shapes will be deactivated.

#### INFO:

**Dynamic rigidbodies** are directly affected by the PhysX solver. Collisions will cause them to move/rotate in response.

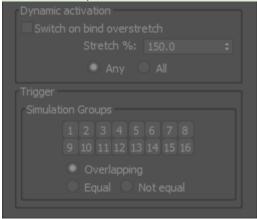
**Kinematic rigidbodies** are not directly affected by the PhysX solver. Dynamic rigidbodies will collide with them, but they themselves will not move in response.

**Trigger rigidbodies** are not directly affected by the PhysX solver. Dynamic rigidbodies will register collision points with triggers rigidbodies, but dynamic rigidbodies with a matching trigger simulation group will not be affected by those collisions.

**Deactivated rigidbodies** are not processed by the PhysX solver at all.

### TIP:

Triggers can be use to trigger PhysX Collision operator test conditions, without causing an actual rigidbody collision response.



## **Dynamic activation**

**Switch on bind overstretch**: causes particles to activate only if attached PhysX bindings stretch past a certain point.

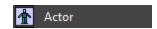
**Stretch** %: the threshold stretch percentage that will cause attached particles to activate.

**Any**: activates particles if any of their PhysX bindings stretch past the stretch threshold.

**All**: only activates particles if all of their PhysX bindings stretch past the stretch threshold.

**Simulation groups**: controls which particle simulation groups will be treated as triggers.

# **Actor operator**



The Actor operator converts **tyActor** objects into hierarchies of actor particles and actor rig particles.



#### Actor

**Input object list**: the list of input **tyActor** objects, and their distribution properties.

#### TIP:

Select an existing item in the actor list to modify its properties.

#### **Distribution**

**Frequency**: the distribution frequency of a particular actor in the list.

#### INFO:

As particles enter the operator, the actor they are assigned depends on the frequency values of shapes in the list. For example, if one actor has a frequency of 10%, and a second actor has a frequency of 90%, there will be a 10% chance that particles will be assigned the first actor, and a 90% chance that particles will be assigned the second actor.

You do not need to ensure that all actor frequencies add up to 100%. Their values will be normalized automatically if they do not all add up to 100%.

## Size

**Override scale**: controls whether to override the scale of particles which are assigned the actor.

Scale %: the scale override.

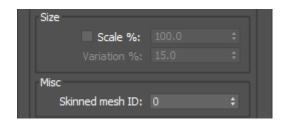
**Variation** %: the per-particle percentage of variation to apply.

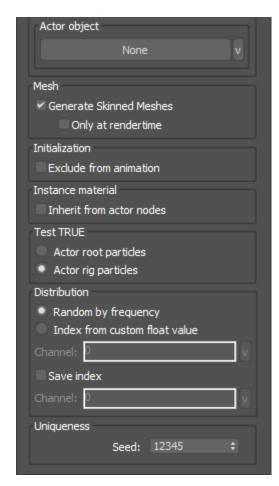
Misc

**Skinned mesh ID**: sets the actor's skinned mesh ID value.

## TIP:

The actor skinned mesh ID value can be used to determine which actor skinned meshes will be exported by an Export Particles operator in tyCache export mode.





### **Actor Object**

**Input tyActor object**: the **tyActor** object that will be used to initialize the actor.

#### Mesh

**Generate skinned meshes**: controls whether **tyFlow** will generate skinned meshes corresponding to input **tyActor** objects which contain skin modifiers.

**At rendertime only**: when enabled, actor skinned meshes will only be generated at rendertime.

#### **Initialization**

**Exclude from animation**: flags actor rig particles, so that they will be excluded by their parent actor particle's animation playback.

#### **Instance material**

**Inherit from actor nodes**: particle render instance materials will be derived from the respective actor nodes.

#### **Test TRUE**

**Actor root particles**: actor root particles will satisfy the test condition.

**Actor rig particles**: actor rig particles will satisfy the test condition.

## Distribution

**Random by frequency**: each particle will choose an actor from the list randomly, taking into consideration each list entry's frequency parameter.

**Index from custom float value**: each particle will choose an actor from the list, using a custom data float value as an index to the listbox's entry array.

Channel: the custom data float channel to retrieve the index value.

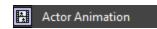
**Save index**: the chosen actor index for each particle will be saved to a custom data float channel.

**Channel**: the custom data float channel to save the index value.

### Uniqueness

**Seed**: the seed value for all varied parameters.

# **Actor Animation operator**



The Actor Animation operator transforms actor rig particles to match pre-defined animation sequences defined in the actor particle's corresponding **tyActor** object.



## **Operation**

**Play animation**: plays a pre-defined animation sequences on any actor particle's applicable actor rig particles.

**Stop animation**: stops the playback of any animation clips on any actor particle's applicable actor rig particles.

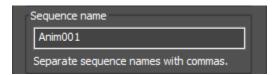
**Include in animation**: flags any previously excluded actor rig particles, so that they will be affected by their parent actor particle's animation playback.

**Exclude from animation**: flags any actor rig particles, so that they will be excluded by their parent actor particle's animation playback.

**Set child coordinates: PARENT**: transforms child rig particles in parent space.

**Set child coordinates: WORLD**: transforms child rig particles in world space.

**NOTE:** By default, **tyFlow** updates rig particles by considering their relation in the corresponding **tyActor** objects they were derived from. If an object added to the **tyActor** is a child of another object added to the **tyActor**, **tyFlow** will take the local transform of the child and update the corresponding rig particle by multiplying that local transform by the particle transform of the equivalent parent rig particle. This way the hierarchy of rig particles is updated in the same manner that the hierarchy of corresponding objects is updated. However, in some situations, it is beneficial to update rig particles independent of their relative parent particle transforms (for example, if the child is a kinematic PhysX particle and the parent is a dynamic PhysX particle and they are bound together with a PhysX bind). In this situation, you can change the relative coordinate system to "WORLD" instead of "PARENT" in order to decouple their transform update.



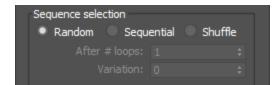
## Sequence name

**Name**: the name of the animation sequences to play on any actor particle's applicNOable actor rig particles.

#### TIP:

To define multiple sequences that will be selected at random for each particle, separate names with commas.

**NOTE:** The names of the sequences must exist in the corresponding **tyActor** object animation sequence list.



## **Sequence selection**

**Random**: the animation sequence for the actor will be chosen at random from the list, when the particle enters the operator.

**Sequential**: all animation sequences listed by the user will play through sequentially, one after the other.

**Shuffle**: a new animation sequence will be chosen from the list, each time the maximum number of loops is reached for the prior sequence.

## **INFO:**

In order for sequential sequences or sequence shuffling to work, operator timing must be set to "Continuous", or a frame/event range that overlaps points in the simulation when shuffling would be expected to occur (ie, when one clip finishes playing).

**After # loops**: controls how many loops to play the current sequence before selecting a new random sequence when shuffling is enabled.

**Variation**: the per-particle amount of variation to apply.



**Blend frames**: the number of blending frames to use when transitioning to a new sequence.

**Blend previous animation**: blending will occur from the previous sequence to the new sequence.

**Blend world transforms**: blending will occur from the previous particle transforms, to the new sequence.

#### **Animation frame range**

**Starting offset**: define a custom starting offset for the sequence.

**Variation**: the per-particle amount of variation to apply.

**Loop**: controls whether sequence playback will loop.

## **Animation playback speed**

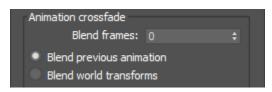
**Speed**: the playback speed multiplier.

**Variation** %: the per-particle percentage of variation to apply.

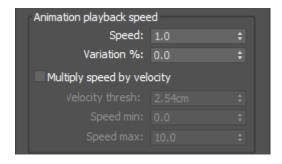
**Multiply speed by velocity**: controls whether playback speed will be relative to particle velocity.

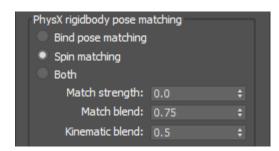
**Velocity thresh**: the velocity threshold which individual particle velocity magnitudes will be relative to.

**Speed min**: the minimum allowed playback speed, relative to the ratio between the velocity threshold and individual particle velocity magnitudes.









**Speed max**: the maximum allowed playback speed, relative to the ratio between the velocity threshold and individual particle velocity magnitudes.

## PhysX rigidbody pose matching

**Bind pose matching**: bindings between particles will be directly adjusted to match relative orientations of input **tyActor** objects.

**Spin matching**: spin forces will be applied to particles in an attempt to match their orientations with the corresponding parts of the input **tyActor** rig.

**Both**: both binding adjustments and spin forces will be applied.

**Match strength**: controls the overall strength of the chosen matching algorithm. A small value will give particles more flexibility in their dynamic motion. A higher value will cause the algorithm to match poses more forcefully.

**Match blend**: the rate at which previously-kinematic rig particles, or rig particles whose actor root particle is new in this operator, will blend into their match pose/spin.

#### TIP:

Reducing the match blend value will allow dynamic PhysX particles to transition into their target poses in a smoother manner. For sufficiently smooth blending, you may need to set this value to be very low (0.01 or lower).

**Kinematic blend**: the rate at which previously-dynamic particles will blend into their kinematic poses.

#### TIP:

Reducing the kinematic blend value will allow kinematic PhysX particles to transition into their target poses in a smoother manner.

Uniqueness



# **Actor Center operator**

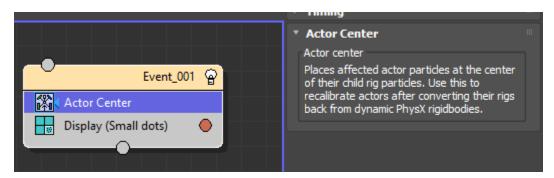


The Actor Center operator places affected actor particles at the center of their child rig particles.

## TIP:

Use this operator to recalibrate actors after converting their rigs back from dynamic PhysX rigidbodies.

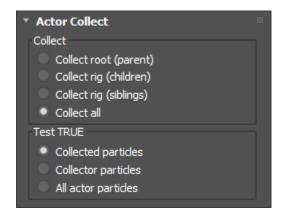
There are no Parameters for this Operator



## **Actor Collect operator**



The Actor Collect operator allows you to collect elements of an actor that have been spread out into separate events, back into the current event.



#### **Collect**

**Collect root (parent)**: brings the parent actor particle into this event

**Collect rig (children)**: brings the children actor rig particles into this event.

**Collect rig (siblings)**: brings the sibling actor rig particles into this event.

**Collect all:** brings all parent/children/sibling actor and actor rig particles into this event.

#### **Test TRUE**

**Collected particles**: collected particles will satisfy the test condition.

**Collector particles**: collector particles will satisfy the test condition.

**All actor particles**: all actor particles will satisfy the test condition.

## **Actor Convert operator**



The Actor Convert operator allows you to convert the skinned meshes of actor particles into particle shape meshes.

There are no parameters for this operator



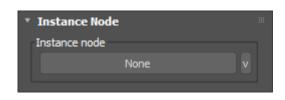
# **Instance Node operator**

The Instance Node operator can be used to assign a render-only node to particles.

#### NOTE:

This is an operator not currently supported by any renderers, so it essentially has no functionality. The reason for its addition is so that renderers can support it in the future, enabling users to scatter any arbitrary objects with particles, not just meshes. In the meantime, users can use the Export Particles (Objects mode) "auto-export on render" option for equivalent functionality.





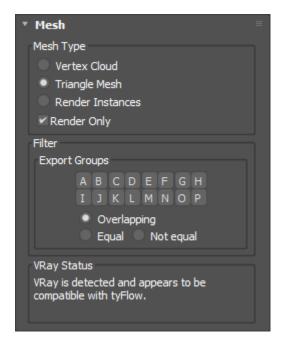
**Instance Node** 

**Node**: the instance node to assign to particles.

## Mesh operator



The Mesh operator can be used to enable particle mesh rendering.



#### INFO:

If you wish to render particles within an event, you must add a Mesh operator to the event.

## **Mesh Type**

**Vertex Cloud**: particles will be converted into a point cloud of isolated mesh vertices.

### TIP:

Vertex Clouds can be rendered by point renderers like Thinkbox's Krakatoa, or converted to isosurfaces using meshing plugins like Thinkbox's Frost.

**Triangle Mesh**: particle shape meshes will be combined into a single editable mesh.

**Renders Instances**: particle shape meshes will be converted into renders instances for compatible renderers.

**Render Only**: controls whether the meshing operation will only happen during rendering.

#### Filter

**Export groups**: controls which particle export groups will be meshed.

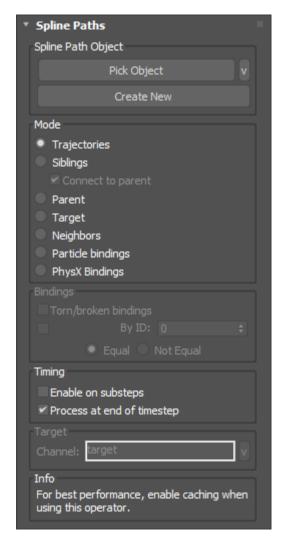
## **Spline Paths operator**



The Spline Paths operator can be used to convert particle trajectories, bindings and relationships into splines.

#### **WARNING:**

Multiple operators within the same flow can send particles to the same **tySplines** object, but you should not send particles from multiple flow objects into the same **tySplines** object, or else the resulting splines will not display correctly.



#### **Spline Path Object**

**tySplines object**: the **tySplines** object that the particle data will be sent to.

#### Mode

**Trajectories**: particle positions will be sent to the **tySplines** object at each step of the simulation, as trajectory data.

**Siblings**: particle sibling data will be sent to the **tySplines** object.

**Connect to parent**: when enabled, the youngest child of a set of siblings will be connected to the parent.

**Parent**: when enabled, particles will be connected to their parent (if they have one).

**Target**: when enabled, particles will be connected to their target (if they have one).

**Neighbors**: particle neighbor proximity data will be sent to the **tySplines** object.

**Particle bindings**: particle binding data will be sent to the **tySplines** object.

**PhysX bindings**: PhysX binding data will be sent to the **tySplines** object.

#### **Bindings**

**Torn/broken bindings**: controls whether data for torn/broken bindings will be included.

**ID enable**: controls whether only bindings matching a certain ID will be included.

**ID**: the ID to match.

**Equal**: bindings will be considered a match if they have the same ID.

**Not Equal**: bindings will be considered a match if they do not have the same ID.

#### **Timing**

**Enable on substeps**: controls whether data will be sent to the input **tySpline** object on simulation substeps.

## Target\*\*

**Channel**: the target custom float data channel.

#### TIP:

"Enable on substeps" can usually be left disabled. It is only necessary when particles are moving so fast that they require extra substep samples for things like motion blur. Enabling this setting will increase the amount of RAM required to store **tySpline** data, relative to the number of substeps required for the simulation.

## **Collision operator**



The Collision operator can be used to collide particles with scene geometry.

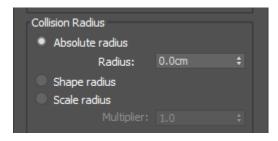
### **Collisions Rollout**



**Input object list**: the list of input collider objects.

#### NOTE:

The Collider operator supports static/deforming geometry objects, as well as built-in Deflector/SDeflector spacewarps.



#### **Collision Radius**

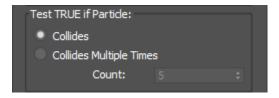
**Absolute radius**: the collision radius of each particle will be set to a specific value.

**Radius**: the specific collision radius value.

**Shape radius**: the collision radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the collision radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

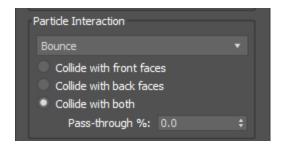


## **Test TRUE if Particle**

**Collides**: the test condition will be satisfied as soon as a particle collides with a collider object.

**Collides multiple times**: the test condition will be satisfied as soon as a particle collides with a collider a specified number of times.

**Count**: the number of times a particle must hit a collider in order to satisfy the test condition.



## **Particle Interaction**

**Bounce**: Particles will bounce off colliders.

**Continue**: Particle will not be affected by colliders (but may still satisfy the test condition).

**Collide with front faces**: particles will only collide with the fronts of faces.

**Collide with back faces**: particles will only collide with the backs of faces.

**Collide with both**: particles will collide with fronts and backs of faces.

**Pass-Through %**: controls the number of particles that are allowed to pass through the collider.

**Geometry Surface Interaction** 

#### Bounce

**Bounce** %: the bounce velocity of particles, relative to their travel velocity.

**Variation %**: the per-particle percentage of variation to apply.

**Diverge %:** the percentage of divergence to apply to bounce vectors.

**Threshold**: particles whose travel velocity magnitude is below this value will not bounce.

#### **Friction**

**Friction** %: the friction applied to particles, relative to their travel velocity.

**Variation** %: the per-particle percentage of variation to apply.

**Threshold**: particles whose travel velocity magnitude is below this value will be treated with 100% friction.

#### **Motion Inheritance**

**Inherit** %: the percentage of collider surface velocity that particles will inherit.

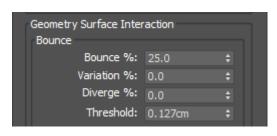
**Variation** %: the per-particle percentage of variation to apply.

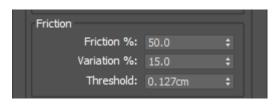
**Apply to Spacewarps**: controls whether Collision operator values relating to bounce/friction/inheritance will override the corresponding values within an input deflector/sdeflector's own settings.

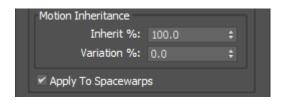
## **Collision Offset**

**Distance**: the distance above a face a particle will be placed after colliding with it.

**NOTE:** Values above 0 help prevent floating-point precision errors during collision detection calculations. Values that are too large will create visible jittering artifact











## **Collision Spawn**

**Mark for collision spawn**: collided particles will be internally marked so that they may satisfy the entry test (collision) condition of a Spawn operator.

**NOTE:** Use this setting in combination with a Spawn operator to spawn child particles at each collision point.

**Evaluation Timing** 

**In order**: the operator will evaluate in order with its surrounded operators.

**After secondary solvers**: the operator will evaluate at the end of the simulation step, after secondary solvers (bind, wobble, etc) have computed their velocities.

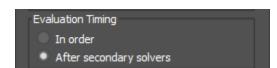
### TIP:

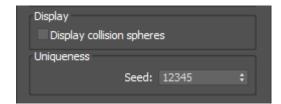
For accurate collisions, the operator should evaluate after secondary solvers (since those solvers may inject velocities into the simulation which will affect collisions). However, in certain circumstances you may not want the operator to evaluate after secondary solvers (if you're using the operator to immediately send particles to another event). In that case, "in order" should be selected as the timing method.

**Display** 

**Display collision spheres**: displays collision spheres for each particle in the viewport.

#### Uniqueness





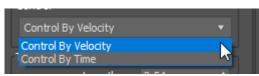
## **Find Target operator**



The Find Target operator provides a way to direct particles towards a specific target point.



#### **Control**



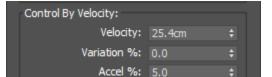
**Control By Velocity**: particles will travel towards their target at a certain velocity.

**Control By Time**: particles will travel to their target within a certain time frame.



**Less than**: particles will satisfy the test condition when their distance to their target is less than this value.

**Variation %**: the per-particle percentage of variation to apply.



Less than: 2.54cm

Variation %: 0.0

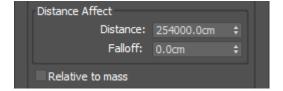
Test TRUE if Distance:

## Control By Velocity\*\*

**Velocity**: the desired travel velocity.

**Variation** %: the per-particle percentage of variation to apply.

**Accel %**: controls how quickly particle forces will accelerate from their old values to the newly applied values.



#### **Distance Affect**

The distance affect extends outwards from the target surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

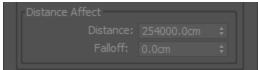


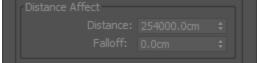
**Relative to mass**: the travel velocity will be relative to the particle's mass.

#### **Control By Time**

**Time**: the time it should take for a particle to reach its target

**Variation**: the per-particle amount of variation to apply.



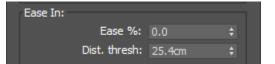


#### **Distance Affect**

The distance affect extends outwards from the target surface.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.



#### Ease In:

**Ease %**: the amount of damping to apply to particle velocities as they approach the target surface.

**Dist. thresh**: the minimum distance to the target surface before the damping takes effect.

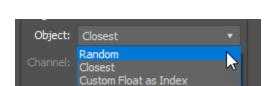
**Target Objects** 

**Input object list**: the list of target input objects.



## **Target Location**

**Object**: the method used to choose a target object from the input object list.



Target Location Object: Closest

**Random**: a random object will be chosen from the list.

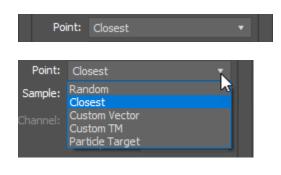
**Closest**: the closest object will be chosen from the list.

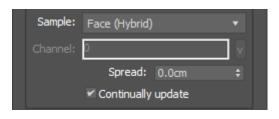
**Custom float as index**: an object will be chosen based on its index in the list. The index value will be taken from a particle's custom float data channel.

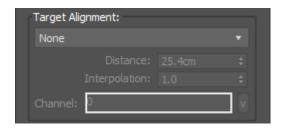


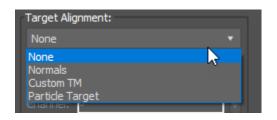
**Channel**: the custom float data channel which will provide the float data.

Cycle through target list: the custom float data value will be incremented, each time a target in the list is reached, until the last target is reached.











**Point**: the method used to choose a target point.

**Random**: a random point will be chosen on the surface of the target object.

**Closest**: the closest point on the surface of the target object will be chosen.

**Custom Vector**: a vector taken from the particle's custom vector data channel will be used as the target point.

**Custom TM**: a vector taken from the particle's custom TM data channel will be used as the target point.

**Particle Target**: a vector taken from the particle's target particle TM will be used as the target point.

**Channel**: the custom data channel which will provide the custom vector data.

**Spread**: the amount of random jitter to add to each target point.

**Continually update**: target points will be updated each step, to account for moving objects and/or changing custom data.

**Target Alignment** 

**Alignment type**: controls how particle orientations are affected by target proximity.

**None**: no changes to particle orientations will be made.

**Normals**: particles will align to target object surface normals as they approach their target point.

**Custom TM**: particles will align to a custom TM data channel value as they approach their target point.

**Particle Target**: particles will align to the TM of their target as they approach their target point.

**Distance**: the minimum distance a particle must be to its target point before its orientation will be interpolated towards its target alignment.

**Interpolation**: the amount to interpolate particle orientations from their previous value to the alignment value.

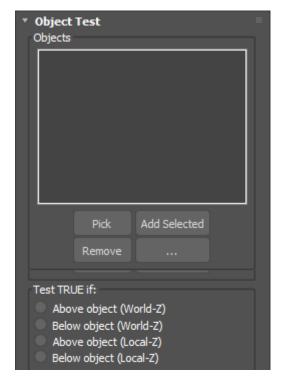
**Channel**: the custom data channel which will provide the custom TM data.

Uniqueness

## **Object Test operator**



The Object Test operator can be used to test particle positions against the positions of objects in the scene.



**Input object list**: the list of input objects to use for the tests.

**NOTE:** Particles will conduct their tests against the closest object in the list of objects.

#### **Test TRUE If**

#### **Above Object (World-Z)**

Particles above the object in world-space along the Z-axis will satisfy the test condition.

#### **Below Object (World-Z)**

Particles below the object in world-space along the Z-axis will satisfy the test condition.

#### Above Object (Local-Z)

Particles above the object in object-space along the object's Z-axis will satisfy the test condition.

### **Below Object (Local-Z)**

Particles below the object in object-space along the object's Z-axis will satisfy the test condition.

**Offset**: the amount of offset that will be applied along the applicable Z-Axis, prior to the test.

**Variation**: the per-particle amount of variation to apply. **Distance to object** 

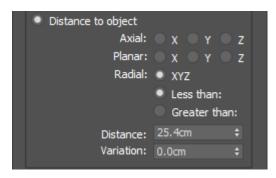
A distance test will be performed between particles and input objects.

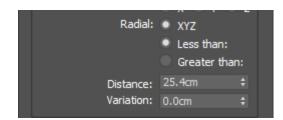
**Axial X/Y/Z**: the distance will be measured between the particle and the object's X/Y/Z axis.

**Planar X/Y/Z**: the distance will be measured between the particle and the plane whose normal is the object's X/Y/Z axis.

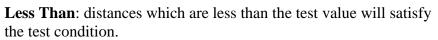
**Radial XYZ**: the distance will be measured between the particle and the object's position.











**Greater Than**: distances which are greater than the test value will satisfy the test condition.

**Distance**: the distance test value.

**Variation**: the per-particle amount of variation to apply.

#### Noise

The noise settings allow you to offset the way in which particle positions are measured during the object test.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

Scale: the scale of the noise.

Roughness: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

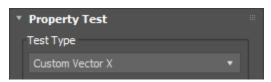
Uniqueness



## **Property Test operator**



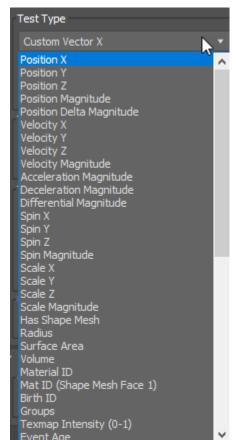
The Property Test operator tests if particles satisfy property-based conditions.

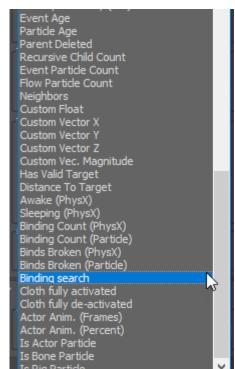


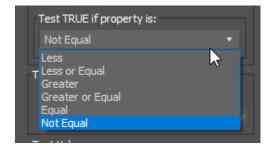
**Test Type** 

**Property List**: the list of testable particle properties.

**Property List**: Screenshot of the list choices of testable particle properties.







**Test TRUE if property is**: various conditions that will be used in the test.

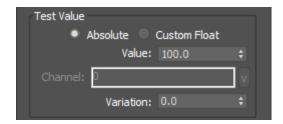


**Test Particle** 

**Self**: the particle's own properties will be tested.

**Target**: properties of the particle's target particle will be tested.

**Channel**: the custom float channel containing the target particle ID.



#### **Test Value**

**Absolute/Custom Float**: controls whether the test value will be taken from the spinner, or a custom float data channel.

Value: the absolute test value.

**Channel**: the custom float channel from which to take the value.

**Variation**: the per-particle amount of variation to apply. **Uniqueness** 

**Seed**: the seed value for all varied parameters.

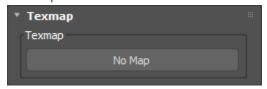


## **Particle Groups Rollout**



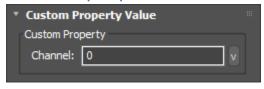
**Simulation groups**: controls which particle simulation groups will be considered for the test.

## **Texmap Rollout**



**Texmap**: the texmap used for the texmap intensity test.

## **Custom Property Value Rollout**



**Channel**: the channel to take the custom float particle data from.

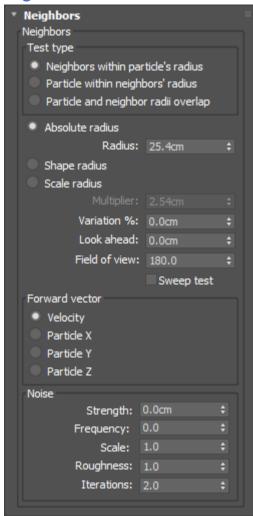
## **Count Binds Rollout**



**Test by ID**: controls whether only bindings which have a matching ID will be counted.

**ID**: the ID to match.

## **Neighbors Rollout**



#### **Test type**

**Neighbors within particle's radius**: a particle will be considered a neighbor if its position is within the searching particle's radius.

**Particle within neighbors' radius**: a particle will be considered a neighbor if the searching particle is within its radius.

**Particle and neighbor radii overlap**: a particle will be considered a neighbor if its radius overlaps with the searching particle's radius.

**Absolute radius**: the neighbor search radius of each particle will be set to a specific value.

Radius: the specific neighbor search radius value.

**Shape radius**: the neighbor search radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the neighbor search radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

**Look ahead**: the search will be conducted at a location this far along the particle's velocity vector.

**Field of view**: the search will be limited to particles within this number of degrees to the particle's trajectory.

**Sweep test**: searches for neighbors along each particle's entire trajectory, not just the particle's starting position.

#### Noise

The noise settings allow you to offset the way in which particle positions are measured during the neighbor test.

**Strength**: the strength of the noise.

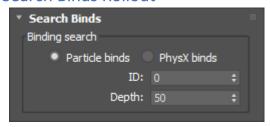
**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

**Scale**: the scale of the noise.

**Roughness**: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

## Search Binds Rollout

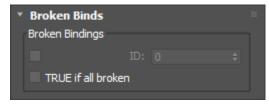


**Particle binds/PhysX binds**: controls which bind type to perform the search on.

**ID**: the binding ID to look for in the bind search.

**Depth**: the maximum depth in the binding hierarchy to search for bindings with the matching ID.

## **Broken Binds Rollout**



**Test by ID**: controls whether only bindings which have a matching ID will be tested.

**ID**: the ID to match.

**TRUE if all broken**: controls whether the test condition will be satisfied if all bindings are broken.

# **Send Out operator**



The Send Out operator can be used to send all particles to the next event

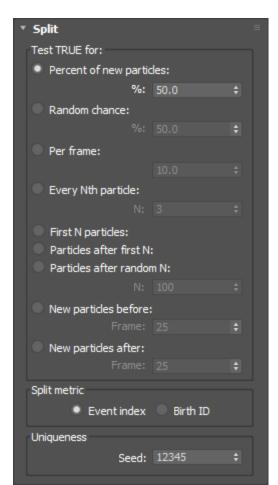
This operator doesn't have any parameters



## **Split operator**



The Split operator tests if particles satisfy conditions based on how they entered the event.



#### **Test TRUE for**

### Percentage of new particles

A percentage of new particles in the event will satisfy the test condition.

**Percentage** %: the percentage value to use.

#### **Random Chance**

Particles will satisfy the test condition based on a random probability.

**Probability** %: the probability value to use.

#### Per frame

A random amount of particles will satisfy the test condition per frame

**Amount**: the amount to send out per frame.

### **Every Nth particle**

Every Nth particle that enters the event will satisfy the test condition.

N value: the N value to use.

#### **First N Particles**

The first Nth particles that enter the event will satisfy the test condition.

**N** value: the N value to use.

#### Particle after first N

Particles after the first Nth particles that enter the event will satisfy the test condition.

**N value:** the N value to use.

#### Particle after random N

Particles after the random N particles in the event will satisfy the test condition.

N value: the N value to use.

### New particles before

Only new particles that enter the event before a certain frame will satisfy the test condition.

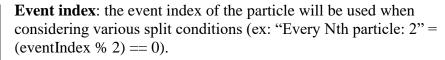
**Frame**: the frame value to use.

## New particles after

Only new particles that enter the event after a certain frame will satisfy the test condition.

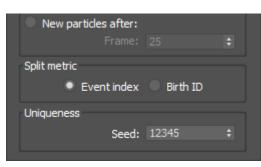
**Frame**: the frame value to use.

## **Split metric**



**Birth ID**: the birth ID of the particle will be used when considering various split conditions (ex: "Every Nth particle: 2" = ((birthID % 2) == 0).

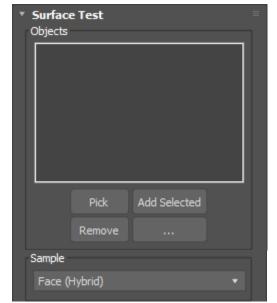
## Uniqueness



# **Surface Test operator**



The Surface Test operator tests if a particle's relationship to a nearby surface satisfies a condition.

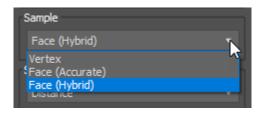


## **Objects**

**Input object list**: the list of input objects to test.

### Sample

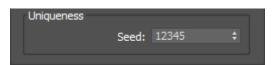
**Sample type**: controls which sampler will be used to determine closest-surface proximities for particles.



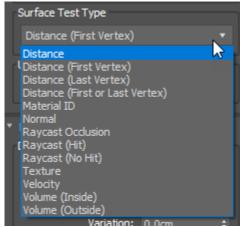


### NOTE:

Depending on the **Surface Test Type** selected from the dropdown menu, various test rollouts appear below the Surface Test Rollout



**Surface Test Type** 



Uniqueness



#### **Inside/Outside Volume**

The test condition will be satisfied if the particle is inside/outside the nearest surface.

**Accuracy**: controls the accuracy of the raycaster used to compute information about whether a particle is inside or outside of an object's volume.



#### Distance to surface

The test condition will be based on the distance between the particle and the nearest surface.



#### Distance to first vertex

The test condition will be based on the distance between the particle and the nearest first vertex.



#### Distance to last vertex

The test condition will be based on the distance between the particle and the nearest last vertex.

#### INFO:

First/last vertex distance tests are useful when trying to determine if a particle following a path has reached the start/end of the path.



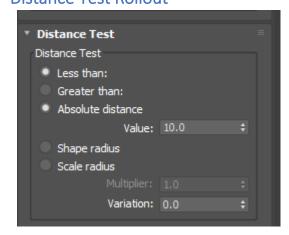
**Distance to edges** 



#### **Raycast occlusion**

The test condition will be satisfied the number of random rays that hit a surface is above a certain threshold (or don't hit the surface if 'invert' is checked)

## Distance Test Rollout



**Less than**: the test condition will be satisfied if the distance is less than a certain value.

**Greater than**: the test condition will be satisfied if the distance is greater than a certain value.

**Absolute distance**: the test distance will be set to a specific value.

**Value**: the specific distance test value.

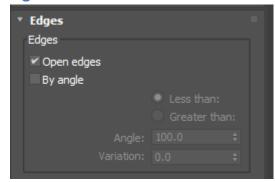
**Shape radius**: the test distance of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the test distance of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.

**Variation**: the per-particle amount of variation to apply.

## **Edges Test Rollout**



## **Edges**

**Open edges**: the test will measure distances to open edges (edges connected to a single face).

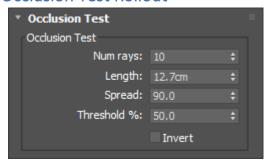
**By angle**: the test will measure distances to edges whose adjacent faces form an angle that matches the specified angle condition.

Less than/greater than: the angle condition.

Value: the angle, in degrees, for the angle condition.

**Variation** %: the per-particle percentage of variation to apply.

## Occlusion Test Rollout



The test condition will be satisfied the number of random rays that hit a surface is above a certain threshold (or don't hit the surface if 'invert' is checked)

**Num rays**: the number of random rays to cast (higher = more accurate).

**Length**: the length of each ray.

**Spread**: the maximum divergence, in degrees, along computed surface normals that the rays will be cast.

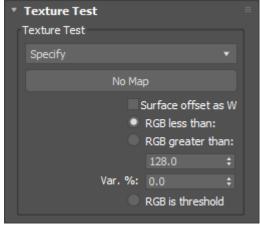
### TIP:

For surface occlusion, a spread of 90.0 is best. For 3d space occlusion, a spread of 180.0 is best.

**Threshold** %: the percentage of rays that must hit a surface to satisfy the test condition.

**Invert**: when enabled, the condition is inverted (it will be satisfied if a percentage of rays above the threshold do *not* hit a surface).

### **Texture Test Rollout**



#### **Nearest Texture**

The test condition will be satisfied if a texture sampled on the nearest point of the nearest surface satisfies a certain condition.

**Texmap**: the texture map to sample.

**Surface offset as W**: the distance from the particle to the closest surface will be used as the W coordinate in the UVW value which is used to sample the texture.

**RGB less than**: the test condition will be satisfied if the magnitude of the sampled RGB value is less than a specified value.

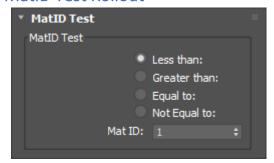
**RGB greater than**: the test condition will be satisfied if the magnitude of the sampled RGB value is greater than a specified value.

Value: the RGB test value.

Variation %: the per-particle percentage of variation to apply.

**RGB** is threshold: the normalized magnitude of the sampled RGB value will be used as test condition probability.

## MatID Test Rollout



#### **Nearest Material ID**

The test condition will be satisfied based on properties of the nearest face's material ID.

**Less than**: the test condition will be satisfied if the nearest face's material ID is less than the specified value.

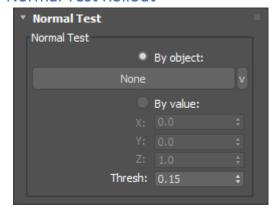
**Greater than**: the test condition will be satisfied if the nearest face's material ID is greater than the specified value.

**Equal to**: the test condition will be satisfied if the nearest face's material ID is equal to the specified value.

**Not Equal to**: the test condition will be satisfied if the nearest face's material ID is not equal to the specified value.

**Mat ID**: the material ID test value.

## **Normal Test Rollout**



#### **Nearest Normal**

The test condition will be satisfied based on properties of the nearest normal on the nearest surface.

**By Object**: tests whether the nearest normal is aligned to the local Z-Axis vector of some object, within the specified threshold.

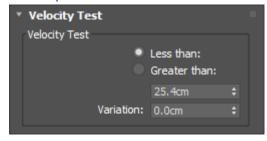
**Object**: the object whose local Z-Axis will be used for the normal test.

**By Value**: tests whether the nearest normal is aligned to a specified normal, within the specified threshold.

**X/Y/Z**: the per-axis values of the specified normal.

**Thresh**: the threshold used to determine normal alignment.

## **Velocity Test Rollout**



#### **Nearest Velocity**

The test condition will be satisfied based on properties of the nearest surface velocity magnitude on the nearest surface.

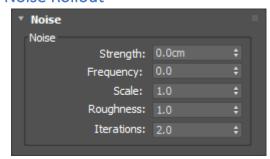
**Less than**: the test condition will be satisfied if the nearest velocity magnitude is less than a certain value.

**Greater than**: the test condition will be satisfied if the nearest velocity magnitude is greater than a certain value.

Value: the magnitude test value.

**Variation**: the per-particle amount of variation to apply.

## **Noise Rollout**



### Noise

The noise settings allow you to offset the way in which particle positions are measured during the distance/volume tests.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

**Scale**: the scale of the noise.

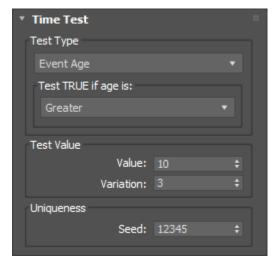
**Roughness**: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

# **Time Test operator**

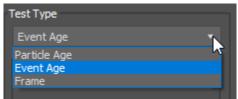


The Time Test operator tests if particles satisfy a time-based condition.



## **Test Type**

**Time type**: the type of time test to conduct.

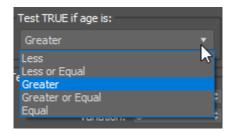


Particle Age: a particle's total age will be tested.

Event Age: a particle's age within its event will be tested.

**Frame**: the current frame of the simulation will be tested.

**Test TRUE if age is**: various conditions that will be used in the test.



### **Test Value**

Value: the test value.

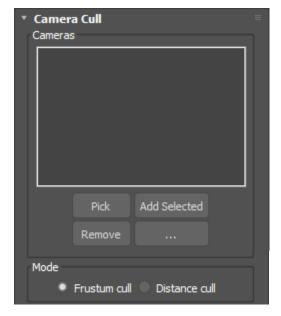
**Variation**: the per-particle amount of variation to apply.

## Uniqueness

## **Camera Cull operator**



The Camera Cull operator can be used to cull particles from display/render/export that are outside of a camera's view frustrum.



#### **Cameras**

**Camera list**: the list of input cameras that will be used to cull particles.

#### Mode

**Frustum cull**: particles will be culled based on whether or not they are inside of a camera's frustum.

**Distance cull**: particles will be culled based on whether or not they are within a certain distance to an object.

**NOTE:** In "distance cull" mode, any object can be used - not just cameras.

Frustum

**Expand**: the number of degrees to expand an input camera's view frustum by.

**Proximity** 

**Distance**: the minimum distance to an input object that a particle must be in order to be culled.

**Invert**: inverts the effect of the distance cull (the distance value becomes a maximum value rather than a minimum value).

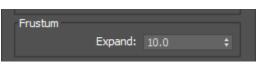
**Affect** 

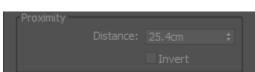
**Display**: when enabled, culling will affect particle display.

**Render**: when enabled, culling will affect particle rendering.

**Export**: when enabled, culling will affect particle exports.

**Particle Interface**: when enabled, culling will affect particle interfaces.







## **Display operator**



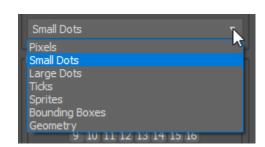
The Display operator can be used to display particles in the viewport.

**NOTE:** Particles are meant to be displayed using the latest Nitrous viewport driver. Displaying particles with a legacy DirectX/OpenGL driver will have a negative impact on performance. Also, not all display types are supported by legacy drivers.



#### **Type**

**Display type**: the drawing method to use for particles.



**Pixels**: particles will be drawn as individual pixels.

**Small dots**: particles will be drawn as small dots.

**Small dots**: particles will be drawn as large dots.

**Ticks**: particles will be drawn as ticks.

**Sprites**: particles will be drawn as images on sprites.

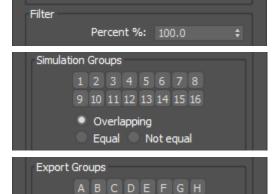
**Bounding Boxes**: particles will be drawn as bounding boxes which encapsulate their shape mesh.

**Geometry**: particles will be drawn as geometry. If a particle does not have an assigned shape mesh, it will be drawn as an X.

#### TIP:

"Sprite" mode is best for granular simulations, where a spherical visualization of particles is necessary. Sprites will display much faster than actual sphere geometry, because they are drawn as simple quads with a sphere texture, rather than actual spherical geometry composed of many vertices and faces.

**Filter** 



IJKLMNOP

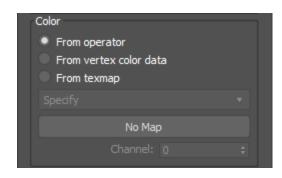
Not equal

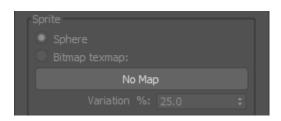
Overlapping Equal N **Percent %**: the percentage of particles to drawn in the view. **Simulation groups** 

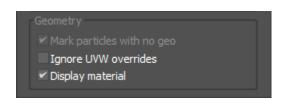
**Simulation groups**: controls which particle simulation groups will be drawn.

### **Export groups**

**Export groups**: controls which particle export groups will be drawn.







#### Color

Color settings apply to non-geometry particle display (ie, all display modes except 'geometry' and 'bounding box'

**From operator**: the color of the particles will be defined by the operator's color picker.

From vertex color data: the color of particles will be defined by the color values assigned to the particle's specified mapping channel.

**From texmap**: the color of particles will be sampled from the specified texmap, using all available mapping channels assigned to the particle.

### **Sprite**

**Sphere**: displays sprites with a built-in sphere texture.

**Bitmap texmap**: displays sprites using a custom bitmap texture map as the texture.

**Variation %**: the amount of color value variation to apply to sprite particles.

#### **Geometry**

**Mark particles with no geo**: particles without a shape mesh will be drawn as an 'X'

**Ignore UVW overrides**: controls whether UVW (mapping) overrides on particles will be ignored.

**Display material**: controls whether geometry will be displayed with the material assigned to the **tyFlow** object. If this is disabled, only the Display operator color will be used, overriding any material assignments.

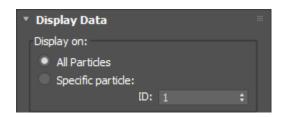
## INFO:

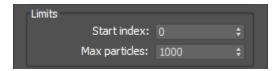
Nitrous GPU instancing does not support UVW (mapping) overrides on instances. Therefore, if particles have active UVW overrides, instead of being instanced they will be combined into a single (potentially gigantic) mesh. Depending on the number of particles in the cache, this can eat up huge amounts of system resources. For example, a million identical particles with no mapping override can be sent to the GPU as a single mesh and a million transforms. But, a million particles with mapping overrides will be sent to the GPU as a million different meshes – something even the most powerful systems will have a hard time processing. By enabling "ignore UVW overrides", mapping overrides on particles will be ignored for viewport display, maximizing the number of particles that can be efficiently instanced in the GPU. The drawback is that with "ignore UVW overrides" enabled, particle mapping channels will not have a visible effect on particle material display in the viewport.

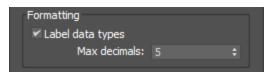
## **Display Data operator**



The Display Data operator allows you to see various particle properties in the viewport.







#### **Display On**

**All Particles**: all particles within the display limits will display their selected data.

**Specific particle**: only the particle with an ID that matches a specific value will have its selected data displayed.

**ID**: the ID to match.

#### Limits

Displaying particle data can be slow, because the Display Data operator uses 3ds Max's legacy text drawing methods. Imposing limits on the total number of particles whose data will be displayed can help to increase performance.

**Start index**: the starting index of particles within the event to display.

**Max Particles**: the number of particles after the starting index whose data will be displayed.

#### INFO:

If the total number of particles in the event is less than (starting index + max particles), the display index will loop back to the beginning of the particle list. So if an event has 10 particles inside it, and the start index is set to 9 and max particles is 5, the operator will draw particles # 9, 10, 1, 2, 3.

**Formatting** 

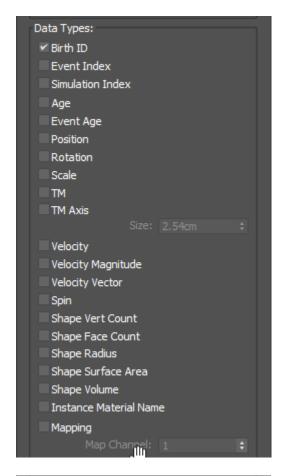
**Label data types**: displays the data type prior to the data value

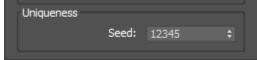
**Max decimals**: the maximum number of decimal places to display for floating point values.

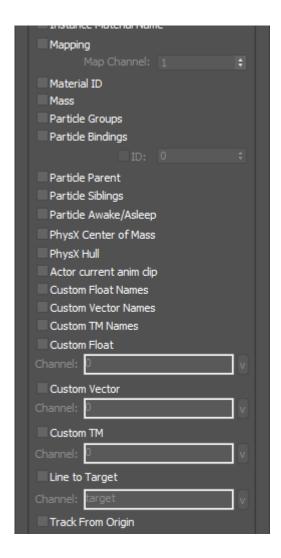
**INFO:** Trailing zeros are automatically removed from floating point values.

#### **Data types**

[**Types**]: the various type of data to draw in the view for each applicable particle.







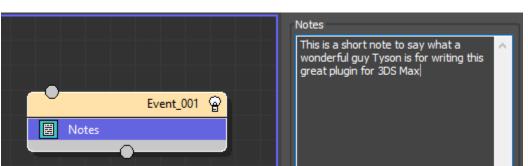
## Uniqueness

# **Notes operator**



The Notes operator allows users to store notes/comments/info inside of a flow.

This operator does not have any parameters, just write whatever notes you want ©



## **Export Particles operator**



The Export Particles operator can be used to convert particles into scene objects or save them into various file formats.

#### INFO:

If an Export Particles operator is placed into an event on its own, or an event only containing other Export Particles operators, its event will operate in "Global Export" mode. This means that the Export Particle operators in that event will operator on all particles within the flow (note: "Global Export" events are shaded blue).

If an Export Particles operator is placed into an event with other regular operators, it operates in "event" export mode. This means it will only export particles in its own event. If an Export Particles operator in "event" mode is instanced across multiple events, it will export particles from all of the events that it's instanced inside.

These modes make it easy to control exactly which particles will be exported by an Export Particles operator.

#### NOTE:

The Export Particles operator will only export particles when its "Export" button is pressed by the user. This operator is not evaluated by the simulation itself.

#### NOTE

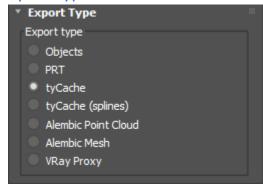
**tyFlow** retimer settings are ignored by the Export Particles operator in PRT or **tyCache** mode. To retime your PRT or **tyCache** sequence, use the retiming spinner in the PRTLoader or **tyCache** object you load your files into. To quickly copy retiming settings between your **tyFlow** object and your cache importer object, right click on the **tyFlow** retiming spinner and choose "Copy Animation", then right click on your cache importer retiming spinner and choose "Paste Animation".

#### **WARNING:**

By default, the Export Particles operator will respect your **tyFlow** cache settings. If caching is enabled, particles that are processed by the Export Particles operator will be added to the **tyFlow** cache. If you attempt to export more particles than can fit in your RAM while caching is enabled, 3ds Max may run out of memory and the export may fail. If you are exporting huge numbers of particles it is recommended to disable the **tyFlow** cache before doing so.

**Note:** This does not apply to export jobs submitted through Deadline, as caching will automatically be ignored by machines that process Deadline tasks.

### **Export Type Rollout**



**Objects**: this mode allows you to convert particles into scene objects.

**PRT**: this mode allows you to convert particles into PRT format files (.prt).

**tyCache**: this mode allows you to convert particles into **tyFlow's tyCache** format files.

**tyCache** (**splines**): this mode allows you to convert splines (generated with a Spline Paths operator, or just regular Max splines) into **tyFlow's tyCache** format files.

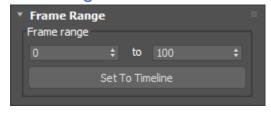
#### **INFO:**

Spline data is stored in a tyCache in a mesh-based format, and can be converted back to proper spline data (post-export) using a **tySplineCache** modifier. Please see the **tySplineCache** modifier documentation for more info.

**Alembic Point Cloud**: this mode allows you to convert particles point clouds into Alembic format files (.abc).

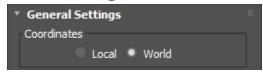
**VRay Proxy**: this mode allows you to convert particles into VRay Proxy format files (.vrmesh).

## Frame Range Rollout



**Start/End**: these spinners control the start/end range of frames over which particles will be exported.

### **General Settings Rollout**



#### **Coordinates**

**Local/World**: the coordinate system that the particles/meshes will be relative to, prior to export.

#### INFO:

When "local" coordinates are chosen, all particle properties will be exported relative to the inverse transform of the **tyFlow** itself. Therefore, the **tyFlow** icon's transform will be the coordinate origin, rather than [0,0,0] in the scene.

**Particles** 



**Export particles**: controls whether particles will be exported. Disabling this setting allows you to limit the export to additional geometry only.

**Birth multiplier**: the birth multiplier applied to birth operators at the time of export. Setting this value to something other than 1.0 will increase/decrease the number of particles generated during export.

**Simulation Groups** 

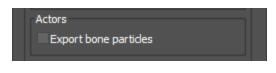
**Simulation groups**: controls which particle simulation groups will be processed by the exporter. Use these groups to limit which particles will be exported.

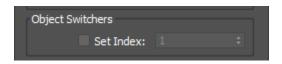
**Export Groups** 

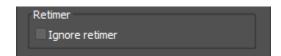
**Export groups**: controls which particle export groups will be processed by the exporter. Use these groups to limit which particles will be exported.











#### **Actors**

**Export bones particles**: controls whether particles imported from a tyActor skin will be exported.

#### **Object switchers**

**Set index**: globally override the switch index of all **tySwitcher** objects in the scene during export.

**Index value**: controls which switch index to use.

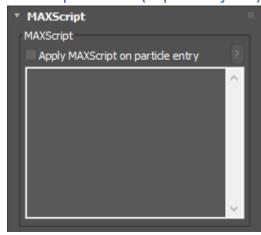
#### TIP:

Use **tySwitcher** objects in your flows to quickly swap between different object setups.

#### Retimer

**Ignore retimer**: turning this on will export raw frame timings (ignoring any retimer settings enabled within the tyFlow).

### **MAXScript Rollout (Export Objects)**



**Apply MAXScript on particle entry**: when enabled, at the time a particle is first exported, custom MAXScript code will be executed on its corresponding scene node.

#### NOTE:

This option allows you to control various parameters of exported objects that would be otherwise too difficult to control from within the **tyFlow** UI. For example, if you are scattering PhoenixFD containers at particle locations, you could use this setting to set their cache frame offset to the time that their particle is born.

### **VRay Proxy Settings Rollout**



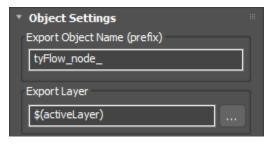
This rollout is exposed when the Export Particles operator is set to "VRay Proxy" mode

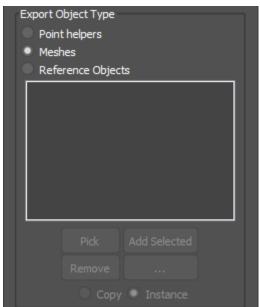
#### **Output**

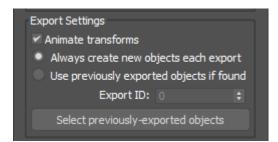
Filename: the output filename for the resulting VRay Proxy file.

### **Object Settings Rollout**

This rollout is exposed when the Export Particles operator is set to "Objects" mode.







#### **Export object name (prefix)**

**Name**: defines the prefix used to name the objects created by the exporter.

**Export layer** 

**Name**: defines the name of the layer that exported objects will be assigned to.

**Export object type** 

**Point helpers**: exports particles as point helper objects.

Meshes: exports particles as editable meshes.

**Reference Objects**: exports particles as duplicates of a reference object taken from the listbox (chosen at random).

**Copy/Instance**: controls whether reference object duplicates will be copies or instances.

#### **Export Settings**

**Animate transforms**: controls whether the transforms of the exported objects will be animated to match the motion of their corresponding particles.

**Always create new objects each export**: each time objects are exported, new scene nodes will be created for all of them.

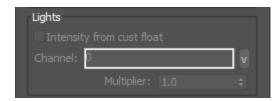
Use previously exported objects if found: if objects were previously exported in the scene, they will be updated during export. Otherwise if such previously-created objects are not found, new ones will be created.

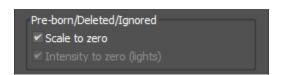
**Export ID**: the export ID is an arbitrary numerical value assigned to exported objects, to help the operator track them.

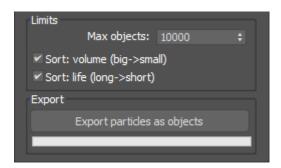
#### INFO:

In order for the operator to track exported objects (for later update), it assigns newly exported objects a tracking code. This code can be viewed in the exported object's user properties. An example code might look something like this:

"tfExport\_10791161940 = 26" and takes the form of "tfExport[operator\_uniqueid][export\_id] = [particle\_id]"







#### Lights

These controls apply to reference objects which are lights (supports standard lights and VRay lights)

**Intensity from cust float**: controls whether the intensity value of the lights will be controlled be particle custom float data.

**Channel**: the custom float data channel.

**Multiplier**: an extra multiplier applied to custom float data values.

Pre-born/deleted/ignored

**Scale to zero**: exported objects that do not have a corresponding particle at a particular frame will have their transform's scale keyframed to zero at that frame.

**Turn off**: exported lights that do not have a corresponding particle at a particular frame will have their intensity animated to zero at that frame.

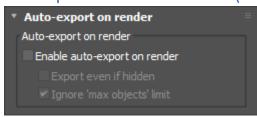
Limits

**Max objects**: controls the maximum number of objects that will be exported.

**Sort: volume**: when the total number of particles exceeds the maximum number of exported objects, the particles will be sorted such that the biggest particles will take priority.

**Sort: life**: when the total number of particles exceeds the maximum number of exported objects, the particles will be sorted such that the oldest particles will take priority.

### Auto-Export On Render Rollout (Export Objects)



**Enable auto-export on render**: when enabled, objects will be exported when a render begins, and subsequently removed when the render ends.

**Export even if hidden**: when enabled, the export will proceed at rendertime even if the **tyFlow** scene object containing the export operator is hidden.

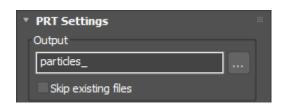
**Ignore 'max objects' limit**: controls whether the "max objects" setting will be respected, while auto-exporting on render.

#### TIP:

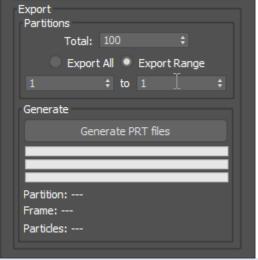
Renderers can typically handle many more objects than the viewport. Since the auto-exported objects will be auto-deleted when rendering ends, you don't have to worry about the viewport slowing to a crawl if you're exporting huge numbers of particles at rendertime.

### **PRT Settings Rollout**

This rollout is exposed when the Export Particles operator is set to "PRT" mode.



#### Channels ✓ ID int32 Age int32 float32[3] ✓ Position float16[4] Rotation float32[3] Scale float32[3] ✓ Velocity Color float16[3] int16 MatID float32 Radius float16[3] TexCoord



#### **Output**

**Filename**: the output filename for the resulting PRT file sequence.

**Skip existing files**: partitions will be skipped if all of their files already exist (and no flow update for that partition will occur). Individual files of an incomplete partition will be skipped if they already exist (but the flow will still need to update for the rest of the partition due to its history-dependent nature).

**Channels** 

[Channel name - data type]: lists the available channels to save in the PRT files, and their corresponding data type.

Data types and their corresponding size in bytes:

byte = 1 byte

int16 = 2 bytes

int32 = 4 bytes

float16 = 2 bytes

float32 = 4 bytes

**NOTE:** The number in square brackets next to some data types represents the number of values that must be stored for that particular channel. For example, a position channel requires X/Y/Z values, each of which is stored as a float32 data type. So the size in bytes for a particular particle position value is 12 bytes (float32 x 3 values).

**Export** 

**Total partitions**: sets the total number of partitions to be demarcated in the PRT filename.

**Export All/Range**: controls whether to export all partitions, or a range of partitions.

#### NOTE:

A PRT file's filename is important, when loading it into a Thinkbox PRTLoader object. Partitions are identified using "partXXofXX" syntax. By keeping the "total partitions" value high, but only exporting a small range of partitions, you can easily add to that range later without creating incompatible filenames. For example, setting "total partitions" to 100 and setting the export range to "1 to 2" will create files marked as

"part01of100" and "part02of100" which will be recognized as two partitions of the same sequence by a PRTLoader. Later you could set the export range to "3 to 3" which would create files marked as "part03of100", etc.

However, if (in that example) you set the total number of partitions to 2, your initial files would be marked as "part01of02" and "part02of02". Later setting the range to "3 to 3" would create files marked "part03of03" which would not be considered part of the same sequence as the prior two partitions, by a PRTLoader.

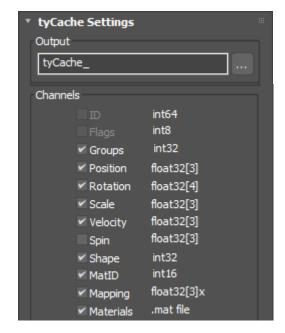
Therefore, it's best to keep the value of "total partitions" high, even if you don't plan on exporting the entire partition range, in order to allow for future increases of the desired partition range within the same sequence.

#### TIP:

By default, all seed values in a flow will be re-seeded during a PRT partition export (for all partitions except for the first partition), in order to randomize particle properties between partitions. If you wish to exclude an operator from the reseeding, simply add the keyword "noseed" to its name. For example, rename "Position Object" to "Position Object noseed".

### tyCache Settings Rollout

This rollout is exposed when the Export Particles operator is set to "tyCache" mode.



#### **Output**

**Filename**: the output filename for the resulting tyCache file sequence.

**Channels** 

[Channel name - data type]: lists the available channels to save in the tyCache files, and their corresponding data type.

Data types and their corresponding size in bytes:

byte = 1 byte

int16 = 2 bytes

int32 = 4 bytes

float16 = 2 bytes

float32 = 4 bytes

NOTE: The number in square brackets next to some data types represents the number of values that must be stored for that particular channel. For example, a position channel requires X/Y/Z values, each of which is stored as a float32 data type. So the size in bytes for a particular particle position value is 12 bytes (float32 x 3 values).

Mapping values are stored as a float32[3] x the number of mapping channels assigned to the particle.



#### Mesh file

**Backup at regular intervals**: controls whether or not the mesh file of the cache (xxx\_tyMesh.tyc) will be updated at regular intervals throughout the export process.

#### INFO:

By default, tyCache mesh files are generated at the end of a tyCache export. If a tyCache export fails (for example: crashes due to lack of available RAM) before the accompanying mesh file is created, you will not be able to load the partial cache. Enabling this option will regularly update the mesh file, so that even if the export fails you may still be able to load the partial cache.

#### NOTE:

For simulations with a lot of unique meshes, the mesh file can get quite large. In those cases, the regular mesh file backups may take a considerable portion of the overall export time. In those cases, it is recommended to disable mesh file backups if you are confident that your export will complete successfully, as then the mesh file export will only need to happen once.

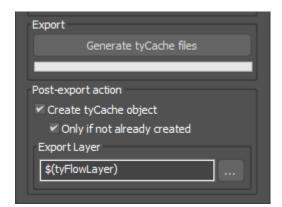
**Post-export action** 

Create tyCache object: controls whether a tyCache object will be created in the scene if the tyCache exporter completes successfully. Its input file sequence will be set to the output sequence of the current exporter, and it will get its material and layer name from the current flow.

Only if not already created: a new tyCache object will only be created, if an existing one using the exporter's output sequence as its input does not already exist. If one already exists, it will simply be updated.

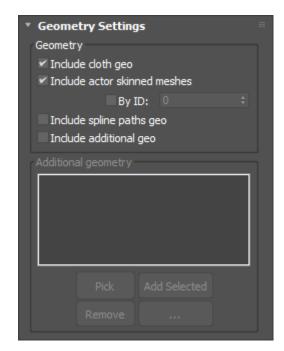
#### **Export layer**

**Name**: defines the name of the layer that created tyCache objects will be assigned to.



### **Geometry Settings Rollout**

This rollout is exposed when the Export Particles operator is set to "tyCache" mode.



#### **Geometry**

**Include cloth geo**: cloth geometry generated with Cloth Bind operators will be added to the **tyCache** on a per-frame basis.

**Include actor skinned meshes**: skinned meshes imported with Actor operators will be added to the **tyCache** on a per-frame basis.

**By ID**: when enabled, only actors with a matching ID will be exported.

**Include spline paths geo**: geometry created with Spline Paths operators (with appropriate tySplineMesher modifiers assigned) will be added to the **tyCache** on a per-frame basis.

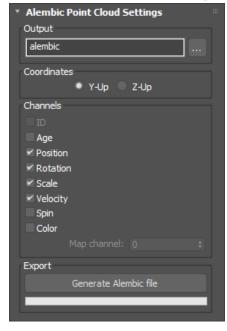
**Include additional geo**: allows you to include additional geometry which will be added to the **tyCache** on a per-frame basis.

#### **Additional geometry**

**Object list**: list of additional objects whose geometry will be included in the **tyCache**.

**NOTE:** An Export Particles operator in "event" export mode that is set to include cloth/actor/spline/additional geometry will include *all* additional geometry (corresponding to those checkboxes) created by the flow - not merely the additional geometry created by its event. This limitation may be subject to change in the future.

### Alembic Point Cloud Settings Rollout



#### **Output**

This rollout is exposed when the Export Particles operator is set to "Alembic Point Cloud" mode.

**Filename**: the output filename for the resulting Alembic file.

#### **Coordinates**

**Y-Up**: exports particles with Y-Up (left-handed) coordinates.

**Z-Up**: exports particles with Z-Up (right-handed) coordinates.

#### TIP:

Y-Up coordinates are used in packages like Maya, Houdini, Unity, etc. Z-Up coordinates are 3ds Max's default coordinates.

#### Channels

[Channel name - data type]: lists the available channels to save in the Alembic file

### **Alembic Mesh Settings Rollout**

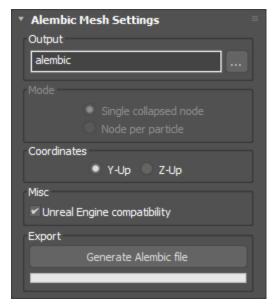
#### INFO:

The Alembic format has evolved over time and not all Alembic importers can properly interpret data exported using the latest Alembic SDK and its various documented methods. Since **tyFlow** uses the latest Alembic SDK, this means that Alembic data exported from **tyFlow** may not be compatible with all importers.

**3ds max:** 3ds Max's Alembic importer uses legacy code that was originally part of the Exocortex plugin suite. In some versions of 3ds Max it is outdated, and does not support important attributes like proper material ID assignments on changing topology. To export Alembic data compatible with 3ds Max's legacy importer, you should export using 3ds Max's own exporter, not the Export Particles operator. Make sure your particles are properly converted to meshes using a Mesh operator before exporting. You should also place a default "Edit Mesh" modifier on your **tyFlow** object so the exporter recognizes it as regular geometry. To import **tyFlow's** exported Alembic data into 3ds Max, you should use an updated importer, like the one included in VRay Proxy objects, which fully supports changing topology. If you use an updated importer, you can export **tyFlow** particles straight from an Export Particles operator. If you use a legacy importer, you need to use the legacy exporter to ensure compatibility.

Maya: Maya's Alembic importer not only relies on legacy Exocortex code too, but it has a buggy implementation which will fail to load **tyFlow** Alembic data or may even crash while attempting to load it. Either export using Max's legacy Alembic exporter (which is compatible with Maya's legacy importer, as long as you follow the legacy export steps listed above), or import into Maya using a VRay Proxy object as the importer.

**Unreal Engine**: Unreal can import **tyFlow's** exported Alembic data so long as it does not include empty frames which contain no geometry (otherwise it may fail to load the data or freeze during playback). For compatibility with Unreal Engine, enable the "Unreal Engine compatibility" checkbox in the Export Particles operator. When that is enabled **tyFlow** will automatically add an infinitesimally small triangle to the cache at all frames, which will prevent Unreal from crashing during playback when no other geometry exists in the cache. When importing into Unreal, choose "geometry cache" mode, set sampling type to "per time step" and set the time step to [1/framerate]. So, for example, for a 30fps sequence, set the time step to 0.03333333.



#### **Output**

This rollout is exposed when the Export Particles operator is set to "Alembic Mesh" mode.

**Filename**: the output filename for the resulting Alembic file.

#### **Coordinates**

**Y-Up**: exports particles with Y-Up (left-handed) coordinates.

**Z-Up**: exports particles with Z-Up (right-handed) coordinates.

### TIP:

Y-Up coordinates are used in packages like Maya, Houdini, Unity, etc. Z-Up coordinates are 3ds Max's default coordinates.

#### Misc

**Unreal Engine compatibility**: since an Alembic cache with no geometry can cause issues in Unreal Engine, this setting will automatically add an infinitesimally small triangle to each frame of the exported Alembic cache, so that Unreal Engine will not find any frames with no geometry in the cache.

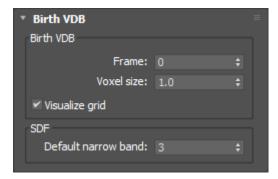
## **Birth VDB operator**



The Birth VDB operator will initialize a new VDB.

#### INFO:

A default VDB is just an empty container. Grids (density, SDF, uvw, etc) must be added to it with the appropriate operators, after it is initialized.



#### **Birth VDB**

Frame: controls which frame the VDB will be initialized on.

**Voxel size**: controls the size of voxels in the VDB.

**Visualize grid**: when enabled, a colored VDB grid marker will be displayed in the view, which also displays the size of the VDB's voxels.

#### **SDF**

**Default narrow band**: controls the default size of SDF narrow bands (the buffer of voxels around the implicit surface of an SDF).

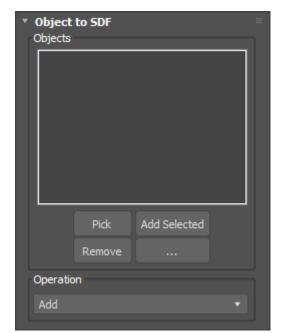
#### TIP:

The default size of 3 is usually best. However, some operations can benefit from a larger narrow band (example: UVW sync during advection on a grid with a tiny voxel size and large velocities. You should usually only increase this value if you see artifacts after advection when using an SDF Solver operator.

# **Object to SDF operator**



The Mesh to SDF operator can be used to convert input meshes into an SDF grid.



#### **Objects**

**Object**: the list of input objects to convert into an SDF grid.

### **Operation**

**Operation**: controls which boolean operation will be used to merge the mesh data into the VDB.

## **Particles to SDF operator**



The Particles to SDF operator can be used to convert input particles into an SDF grid.



#### **Particles**

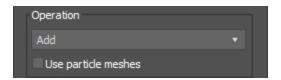
**Include this flow's particles**: when enabled, the particles in the same flow as this operator will be input into the VDB.

**Integrate spin/velocity**: when enabled, the same flow's particles will have their spin/velocity integrated during the input process.

#### NOTE:

Normally spin/velocity integration happens for a particle at the end of the timestep. Therefore, in order to properly sync VDB data with particle data of the current timestep, spin/velocity must be manually integrated beforehand. Hence, the need for this setting.

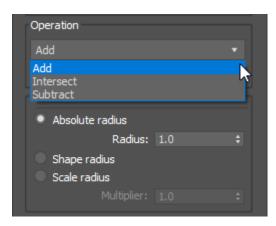
**Particle objects**: the list of additional particle systems to be input into the VDB.



#### **Operation**

**Operation**: controls which boolean operation will be used to merge the particle data into the VDB.

**Use particle meshes**: when enabled, each particle's mesh will be input into the VDB, instead of simply considering the particle as a point with a radius.



#### **Radius**

**NOTE:** When "use particle meshes" is disabled, each particle will be treated as a point with a radius (ie, a sphere).

**Absolute radius**: the radius of each particle will be set to a specific value.

**Radius**: the specific radius value.

**Shape radius**: the radius of each particle will be set to each particle's shape mesh radius.

**Scale radius**: the radius of each particle will be set to each particle's maximum scale dimension.

**Multiplier**: the multiplier to apply to shape/scale radius values.



## **Filtering**

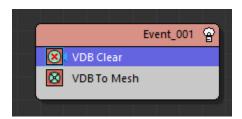
**Simulation/export groups**: controls which groups will be used to filter input particles.

# **VDB** Clear operator



The VDB Clear operator can be used to clear all voxels in a VDB.

This operator has no parameters.



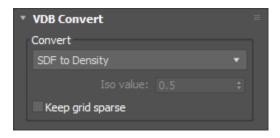
## NOTE:

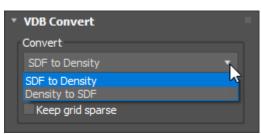
All voxels within all grids of the VDB will be cleared.

## **VDB** Convert operator



The VDB Convert operator can be used to convert SDF voxels to Density voxels and vice versa.





#### Convert

**Mode**: controls which conversion mode (SDF to Density or Density to SDF) will be activated.

**Iso value**: controls the iso value for density grid conversion. Higher values bias the surface level towards higher densities.

**Keep grid sparse**: when enabled, simplified voxel groups in the VDB tree will not be broken up into individual voxels.

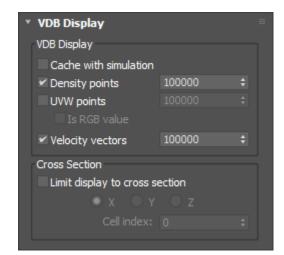
#### NOTE:

"Keep grid sparse" can reduce memory usage, but if voxels values are significantly changed in some way (ex: displaced), large square artifacts can appear in the grid.

## **VDB** Display operator



The VDB Display operator can be used to display VDB data in the viewport.



#### **VDB** Display

**Cache with simulation**: when enabled, display data (vectors, points, etc) will be computed and stored with every frame of the simulation.

#### NOTE:

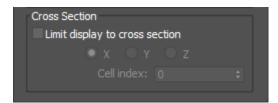
Due to the amount of data contained in a VDB grid (often times, many millions of voxels), storing all relevant display data can be both computationally and memory expensive. Thus, the ability to exclude VDB display data from the **tyFlow** cache has been added. However, since VDB data is not cached on a per-frame basis either, when "cache with simulation" is disabled, only data for the most recent frame of the simulation will be displayed. So if you need per-frame visualization data you can enable this setting, but be mindful that it may eat up a lot of memory and slow down the simulation.

**Density points**: when enabled, a specified maximum number of density voxels will be displayed as grayscale points.

**UVW points**: when enabled, a specified maximum number of UVW voxels will be displayed as colorized points.

**Is RGB Value**: when enabled, the UVW voxel values will be treated as 0-255 RGB values, therefore affecting how they are colored.

**Velocity vectors**: when enabled, a specified maximum number of velocity voxels will be displayed as colorized lines. **Cross Section** 



#### INFO:

In large, complex grids...displaying all relevant grid information in the viewport can be too messy. However, by enabling the cross-section settings, you can view a single slice of the grid at a time.

**Limit display to cross section**: enabled cross section display mode.

**X/Y/Z**: controls which axis of the VDB the cross section will be taken from.

**Cell index**: controls which cell index of the specified axis the cross section will be taken from.

## **VDB** Filter operator

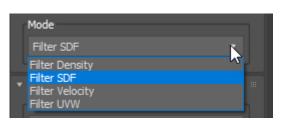


The VDB Filter operator can be used to apply various filtering algorithms to voxels.

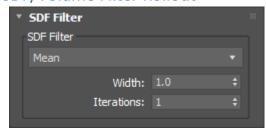
#### Filter Mode Rollout

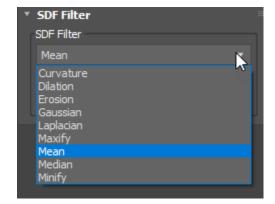


**Mode**: controls which VDB grid will be filtered.



## SDF/Volume Filter Rollout





**Filter**: controls which filter will be applied to the specified grid.

#### NOTE:

Explanations for each filter can be found in the online OpenVDB documentation, with the exception of "minify" and "maxify", which are filters that apply dilation/erosion in equal amounts in order to reduce isolated voxels or fill concave holes in voxels clusters.

Width: the size of the filter (in voxel units).

**Iterations**: the number of times to apply the filter each time step.

## **VDB Modify operator**



The VDB Modify operator can be used to mathematically transform voxels in a VDB.

#### INFO:

The VDB Modify operator is at the heart of all VDB operations within **tyFlow**. It can be used to initialize various grid types (velocity, uvw), composite voxel values together, transfer surface properties, etc. Understanding how this operator works and harnessing its abilities is the key to effectively working with VDBs within **tyFlow**.

#### NOTE:

The documentation for this operator is not yet complete, and will be uploaded at a later date.

### **Modify Grid Rollout**



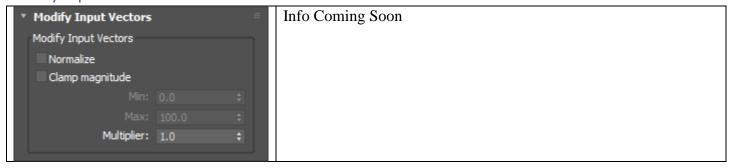
## **Vector Operation Rollout**



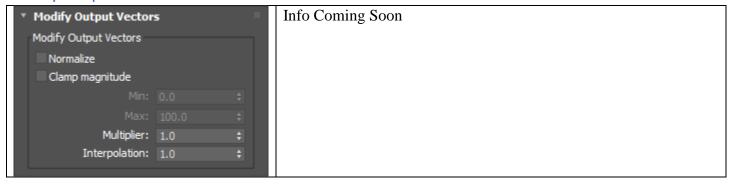
## Input Vectors Rollout



## **Modify Input Vectors Rollout**



## **Modify Output Vectors Rollout**



## Info Rollout



## **VDB Solver operator**



The VDB Solver operator can be used to advect volumes with a velocity field.

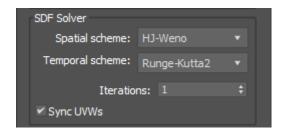


#### Mode

**Mode**: controls which grid type will be advected.



**SDF Solver** 

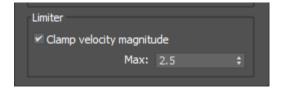


**Spatial/temporal scheme**: controls which scheme will be used to advect and then re-normalize the SDF grid. The default schemes tend to produce the best results in the shortest amount of time.

**Iterations**: the number of iterations of advection to apply to the grid.

**Sync UVWs**: when enabled, any existing UVW grids will be advected in lockstep with the SDF grid, using a custom particle-based proximity scheme.

Limiter



#### NOTE:

Since the advection algorithm traces voxels along their trajectory as they are advected, large velocity can cause significant calculation slowdowns. Enabling the limiter can alleviate this problem, with the caveat that clamping velocities can lead to artifacts if many voxel velocities are above the maximum value. These settings mostly act as a safety, whereas the proper workflow is to simply reduce velocities or reduce simulation timestep.

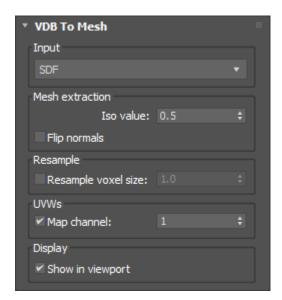
**Clamp velocity magnitude**: when enabled, the magnitude of voxels in the velocity grid used to perform the advection will be clamped.

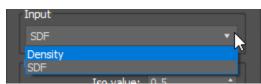
**Max**: the maximum allowable magnitude for any voxel in the velocity grid.

## **VDB** to Mesh operator



The VDB to Mesh operator can be used to convert VDB grids into a mesh.





#### **Input**

**Grid type**: controls which VDB grid will be input into the mesh conversion algorithm.

#### **Mesh extraction**

**Iso value**: controls the iso value for density grid conversion. Higher values bias the surface level towards higher densities.

**Flip normals**: flips the normals of the resulting mesh.

#### Resample

**Resample voxel size**: when enabled, the input grid will be resampled to the specified voxel size prior to mesh conversion.

#### **UVWs**

**Map channel**: when enabled, any existing UVW grids will be applied to the resulting mesh's specified map channel.

#### **Display**

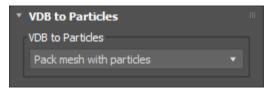
**Show in viewport**: controls whether the resulting mesh will be displayed in the viewport.

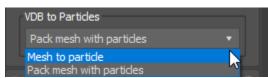
## **VDB** to Particles operator



The VDB to Particles operator can be used to convert VDB meshes into particles.

#### **VDB** to Particles Rollout

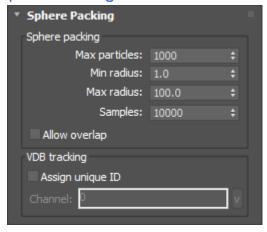




**Mesh to Particle**: converts any generated VDB meshes into a single particle.

**Pack mesh with particles**: packs any generated VDB meshes with particles.

### **Sphere Packing Rollout**



#### Sphere packing

**Max particles**: the maximum number of particles to pack the VDB mesh with.

Min/max radius: radial constraints applied to generated particles.

**Samples**: the number of samples to use in the packing algorithm. Higher values increase packing accuracy.

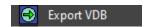
**Allow overlap**: when enabled, resulting particles are allowed to overlap. When disabled, no generated particles will overlap.

#### **VDB** tracking

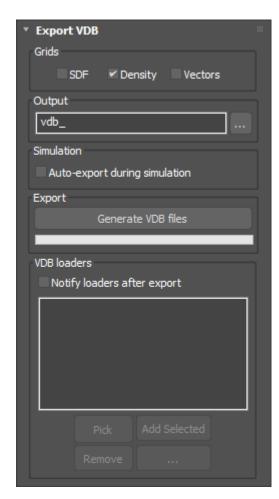
**Assign unique ID**: when enabled, each iteration of the sphere packing algorithm will assign a unique ID to the resulting particles.

**Channel**: the name of the custom float data channel where the unique ID will be assigned.

## **Export VDB operator**



The Export VDB operator can be used to export VDB data to disk.



#### Grids

**SDF/Density/Velocity**: controls which VDB grids will be saved to the .vdb file.

#### **Output**

**Filename**: the output path for the .vdb file.

#### **Simulation**

**Auto-export during simulation**: when enabled, .vdb files will be written automatically as the simulation progresses, after each time step.

#### **WARNING:**

Enabling this setting will cause any existing .vdb files in the output path to be overwritten, any time the simulation progresses.

#### **Export**

**Generate VDB files**: when pressed, the entire timeline will be simulated, exporting .vdb files after each timestep.

#### NOTE:

Because VDB data is not cached on a per-frame basis, a previously-cached sim must be recomputed during manual export.

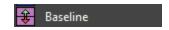
#### **VDB** loaders

**Notify loaders after export**: when enabled, the specified VDB loaders will be refreshed after the .vdb file(s) are exported.

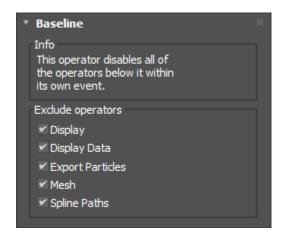
#### **INFO:**

The Export VDB operator will attempt to notify scene VDB loaders that their data needs to be updated, but only some loaders are officially supported for this task. The officially supported loaders are: PhoenixFDFire containers and RedshiftVolumeGrid containers.

## **Baseline operator**



The Baseline operator disables all of the operators below it within its own event.



#### TIP:

Sometimes, in order to step through an event to see how it's affecting particles, you may want to disable and then progressively re-enable operators in the event from top to bottom. This can be a tedious process if your event has a large operator stack. The Baseline operator allows you to quickly disable all other operators below it within an event. The Baseline operator is not allowed to overwrite an existing operator in an event, so you don't have to worry about accidentally dropping it on top of another operator as you move it around.

**Exclude operators**: controls which operator types will be excluded from the disable operation.

# **Debug operator**

The Debug operator prints particle data to a file, for debugging purposes.

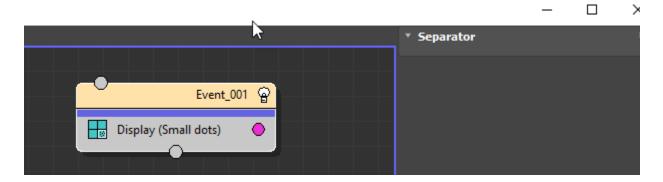
Note: as of 5/6/2020 I didn't see this operator in the actual Max interface, but am leaving it in the documentation for completeness – it is in tyFlows Documentation.

# Separator



The Separator allows you to create a visual separation between operators in an event.

There are no parameters for this operator



## tyFlow Controllers

## tyParticleController Controller

The **tyParticleController** transform controller is a controller that can be assigned to any object's transform, in order to bind that object to a particular particle of a particular flow. This allows you to drive an object's motion directly, using particles.

#### **Particle Controller**

- **tyFlow Object**: choose the **tyFlow** object whose particles will be used to drive the controller.
- **Particle ID**: choose the particle ID from the flow whose transform will by read by this controller, driving its motion.

## tyRateOfChange Controller

The **tyRateOfChange** float controller is a controller that can be assigned to any animatable float value. Its purpose is to provide an easy and procedural way to assign keyless linear animation to a float.

#### INFO:

If you've ever wanted a particular float value to change steadily over time (ex: Noise texmap phase parameter) without having to worry about setting keyframes on it, the **tyRateOfChange** controller is an easy way to do it.

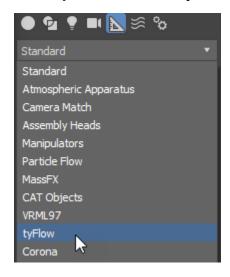
- Rate of change: the rate at which the float value will change over time.
- **Ref. frame**: the reference frame from which changes will begin to occur.
- Value offset: an overall offset that will be applied to the resulting float value.
- **Per frame/second**: the unit of time in which the rate of change will be calculated.
- Clamp value: controls whether the final value will be clamped within a min/max range.
- Min/max: the min/max range in which the final value will be clamped.

# tyFlow Helpers

There are several **tyFlow Helpers**, we will go over each one in the next few pages, they can be accessed from the helpers tab.



Select **tyFlow** from the helper's selection drop down menu

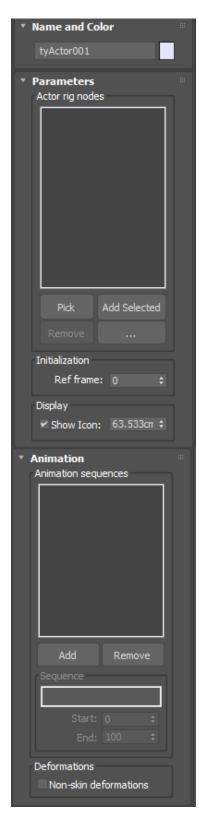


This brings up the helper types to select from



## tyActor Helper

The **tyActor** helper can be used to transfer character rigs into **tyFlow** for use in crowd simulations, in a way that maintains skinning weights, object hierarchies, and keyframed animation of the input objects.



#### **Parameters**

**Actor rig objects**: All objects added to this list will be added to the actor, and processed by **tyFlow**.

**Show icon:** controls whether the **tyActor** icon is displayed in the viewport.

**Icon size**: controls the size of the **tyActor** icon.



#### **Animation**

**Animation sequences**: a list of sequences added to the **tyActor** object, for use in the corresponding flow.

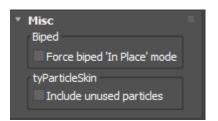
**Sequence name**: takes string input for the name of a particular animation sequence being added.

**Start**: the starting frame of the sequence.

**End**: the ending frame of the sequence.

**Non-skin deformations**: controls whether non-skin deformations are to be considered during sequence playback by the corresponding flow (morpher/bend/twist/etc: anything that transforms vertices over time). Turning this on allows users to assign any vertex animation onto the resulting actor particles, not simply skin deformations.







Triggers are objects that the oriented bounding boxes of the rig objects can intersect with. If an intersection occurs during the animation clip, corresponding rig particles' custom data channel will be modified to hold the 'triggered' value. When no intersection occurs, the rig particles' custom data channel will be modified to hold the 'untriggered' value.

**Triggers**: the list of trigger objects that the **tyActor** will use.

#### **Trigger settings**

**Trigger object**: the scene object used to perform the intersection test against the rig objects.

**Channel**: the name of the custom data float channel to apply the triggered/untriggered values to.

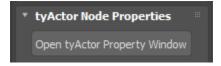
**Untriggered**: the value that will be assigned to corresponding rig particles when no intersection is occurring.

**Triggered**: the value that will be assigned to corresponding rig particles when an intersection is occurring.

Misc

**Force biped "in place" mode**: this is a workaround to solve the problematic "In Place" mode of character studio bipeds, which does not stay enabled in certain circumstances. This setting will force "In Place" mode to be enabled, prior to rig evaluation by **tyFlow**.

**Include unused particles**: by default, particles referenced by a **tyParticleSkin** modifier processed by **tyActors** will be culled if no vertices are weighted to them. Enabling this setting prevents the particle culling from occurring. All particles referenced by a **tyParticleSkin** will be imported, even if no skinned vertices are weighted to them.



### tyActor object properties

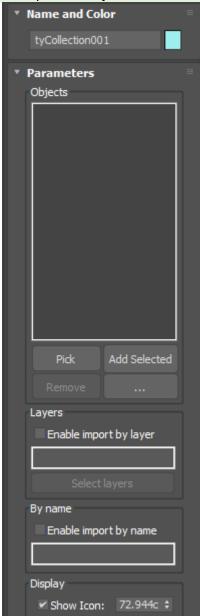
**Open tyActor property window**: opens the floating property window for **tyActor** rig objects. You can use this window to override particle properties of objects (for later use in the corresponding flow). Properties are stored as *User defined properties* on the affected objects.

## tyCollection Helper

The **tyCollection** helper allows you to create a list of objects that can be input into a flow as a single object (the tyCollection object itself), as opposed to adding the same list of objects to multiple operators manually.

#### TIP:

Instead of manually managing large groups of objects shared between multiple **tyFlow** operators, which can be a tedious process, simply input the list of objects into a **tyCollection** and then add the **tyCollection** to the operators instead. When the list of objects in a **tyCollection** is changed, all relevant operators will be sent the updated object list automatically.



**Parameters** 

#### **Objects**

**Objects**: the list of **tyCollection** objects.

#### Layers

**Enable import by layer**: enables the import of objects into the **tyCollection** by layer name.

**Layers**: the names of the layers whose objects will be imported into the **tyCollection**.

#### TIP:

For multiple layer names, separate names with commas (ex: "Layer001, Layer002, Layer003")

#### By name

**Enable import by name**: enables the import of objects into the **tyCollection** by name.

**Names**: the names of the objects which will be imported into the **tyCollection**.

#### TIP:

Asterisks can be used as name wildcards (ex: "Sphere" to import Sphere001, Sphere002, Sphere003, etc). For multiple object names, separate names with commas (ex: "Sphere, Box001, Cylinder00\*")/

#### **Display**

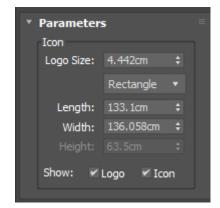
**Show icon:** controls whether the **tyCollection** icon is displayed in the viewport.

**Icon size**: controls the size of the **tyCollection** icon.

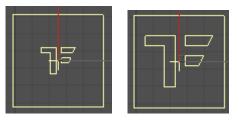


# tyIcon Helper

The **tyIcon** helper can be used by various **tyFlow** operators to control the position or direction of particles.

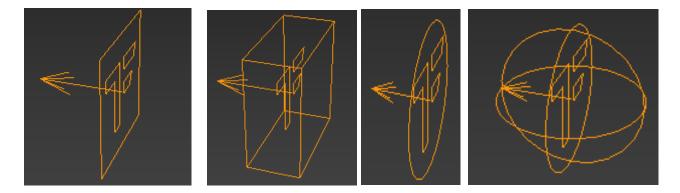


**Logo size**: controls the size of the **tyIcon** logo in the viewport.



**Icon type**: controls the icon shape type.





**Length/Diameter**: controls the length or diameter of the icon, depending on the icon shape type.

Width: controls the width of the icon, depending on the icon shape type.

**Height**: controls the height of the icon, depending on the icon shape type.

**Show logo/icon**: controls which aspect of the icon are drawn in the viewport.

## tyProp Helper

The **tyProp** helper can be used to randomize actor props within a crowd simulation. By adding a **tyProp** to your **tyActor** object (and attaching it to your input character's hierarchy, in place of the actual prop you want to randomize), **tyFlow** will choose a random object from its prop list for each instance of the actor in the flow, effectively randomizing props between multiple actors.

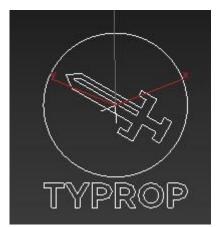


#### **Parameters**

**Prop list**: the list of input props which will be assigned, at random, to instances of the parent **tyActor** object, in place of the **tyProp** object.

**Show icon**: controls whether the **tyProp** icon is displayed in the viewport.

**Icon size**: controls the size of the **tyProp** icon.



# tySlicer Helper

The **tySlicer** helper can be used to manually slice particle bindings. By treating it like a virtual knife, you can animate it slicing across particle bindings within your scene and the corresponding bindings will break.



#### **Parameters**

**Length**: the length of the slicer. The "blade" of the slicer must cross over a binding in order to break it, so the length of the slicer will affect whether such an intersection takes place.

**Show logo**: controls whether the **tySlicer** logo is displayed in the viewport.



## tySwitcher Helper

The **tySwitcher** helper allows you to quickly iterate between different input objects in a flow. For example, sometimes you may have a complex flow setup that you want to transfer between different objects in a scene (with the only difference between setups being their input objects). In that situation you would normally just duplicate your **tyFlow** object and re-assign the alternative objects to the relevant operators in the duplicated flow, keeping all other aspects of the flow the same besides those differing inputs. The **tySwitcher** is an easy way to go about doing that much more efficiently.

By adding a list of input objects to a **tySwitcher**, you would simply choose the **tySwitcher** *itself* as the input object for the relevant operators in your flow, in any place where an input object is accepted (position icon, surface test, etc). Then, by modifying the switch index of the **tySwitcher**, you can choose which object in the list the **tySwitcher** will pass up to the flow. The flow operators will take the input object sent to it by its switchers, and simulate its particles using those objects.



#### **Parameters**

**Objects**: the list of switcher objects.

**Switch index**: the index of the object in the object list that will be passed to any flow operators that the switcher has been added to.

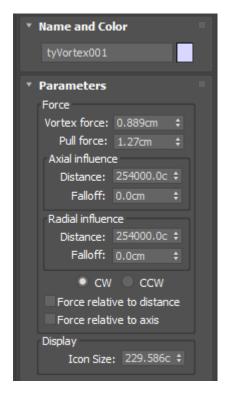
**Show icon**: controls whether the **tySwitcher** icon is displayed in the viewport.

**Icon size**: controls the size of the **tySwitcher** icon.



# tyVortex Helper

The **tyVortex** helper can be used as an alternative to 3ds Max's *Vortex* spacewarp, within a flow. It applies circular forces to particles along an axis.



#### **Parameters**

**Vortex force**: controls the vortex force applied to particles, tangential to the vortex central axis.

**Pull force**: controls the pull force applied to particles, which pulls particles towards its central axis.

# **Axial influence**

The axial influence extends outwards from the vortex's vertical axis.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

#### Radial influence

The radial influence extends outwards from the vortex' center.

**Distance**: particles within this distance will be fully affected.

**Falloff**: the effect on particles beyond the base distance, but within this falloff distance, will diminish according to the inverse-square law.

**CW/CCW**: controls the direction of the vortex (clockwise or counterclockwise).

**Force relative to distance**: controls whether the force applied to particles will be relative to their distance to the vortex.

**Force relative to axis**: controls whether forces applied to particles will be constrained to ensure that particle trajectories remain tangential to the central axis of the vortex. Turning this on will help ensure particles move in a perfect circle around the central axis regardless of their velocity, rather than being flung away, especially when "Force relative to distance" is turned on.

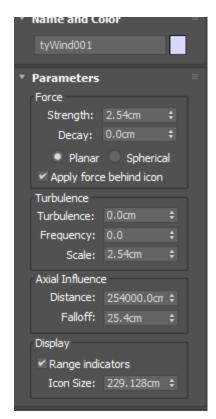
**Icon size**: controls the size of the **tyVortex** icon.



# tyWind Helper

The **tyWind** helper can be used as an alternative to 3ds Max's *Wind* spacewarp, within a flow. It has extra axial falloff controls.

# tyWind Rollout



#### **Force**

**Strength**: controls the overall strength of the wind force.

**Decay**: controls the distance decay of the wind force.

**Planer/Spherical**: controls the outward direction of the wind force. Planar wind forces point in the direction of the icon arrow. Spherical wind forces point in all directions out from the center of the wind icon.

**Apply force behind icon**: when enabled, particles below the **tyWind** force plane's local z-axis will be affected by the **tyWind** forces. When disabled, they will not be affected.

## **Turbulence**

**Turbulence**: the strength of the wind's turbulence.

**Frequency**: the frequency of the wind's turbulence.

**Scale**: the scale of the wind's turbulence.

# **Axial Influence**

**Distance**: particles within this distance to the z-axis of the wind will be fully affected by the wind.

**Falloff**: the effect on particles beyond the base distance to the z-axis, but within this falloff distance, will diminish according to the inverse-square law.

# **Display**

**Range indicators**: draws helpers shapes which represent the decay and axial falloff values in the viewport.

**Icon size**: controls the size of the **tyWind** icon.

# **Colliders Rollout**

Adding meshes to the colliders rollout allows you to occlude particles from the **tyWind** force, if a colliding face separates them from the origin of the wind force on the **tyWind** helper.



#### Collider list

**Mesh list**: the list of meshes that may block wind forces from affecting particles.

# **Collider affect**

**Ray divergence**: controls how many degrees to diverge force rays cast from the **tyWind** object towards particles. The larger the value, the less sharp blocking edges around colliders will be.

Occlusion %: the amount to occlude particles from forces if a collider is hit.

# TIP:

Lower this value to simulate semi-occluding materials (ex: a screen door or thin cloth material). This value is additive across multiple colliders that are hit, so that if you set it to less than 100%, the more colliders that are hit, the more force-occlusion a particle will undergo.

# tyFlow interaction

**Collide with cloth meshes**: when enabled, cloth meshes of the source tyFlow will be added to the list of mesh colliders.

# tyMaterials

The **tyMtlLib** material automatically loads a material from a .mat file, for simple material referencing. Changes to a .mat file will automatically propagate to any **tyMtlLib** materials, when a max file is reloaded, or the "reload" button is pressed in the **tyMtlLib** UI. The **tyMtlLib** functionality is similar to that of the built-in XRefMaterial, however loading/referencing materials directly from a .mat can be more convenient than loading/referencing them from an XRef.

# tyLibMtl

• **Filename**: the name of the .mat file to load.

# **Material selection**

• **Index**: the 0-based index of the material in the .mat file to use.

# NOTE:

The chosen index will be internally clamped to the range of available materials in the .mat file.

# **MAXScript**

# **MAXScript Editor Access**

**tyFlow** objects have MAXScript interfaces which can be used to manipulate their editor windows.

- [obj].editor\_open(): opens the tyFlow editor GUI.
- [obj].reset\_simulation(): resets the tyFlow simulation and clears the cache.

# **MAXScript Particle Access**

Both **tyFlow** and **tyCache** objects have MAXScript interfaces which can be used to read particle data.

# **Preparing particles**

• [obj].updateParticles {frame}: prepares particle data for MAXScript access at a particular frame.

#### **WARNING:**

The **updateParticles** function of a **tyFlow** or **tyCache** object **MUST** be called prior to accessing particle data.

• **[obj].numParticles**(): returns the number of particles at the prepared frame.

#### **Individual particle data**

Data for individual particles can be accessed with their index. Valid particle indices (for **tyFlow/tyCache** objects with at least 1 particle) range from **1** to **[obj].numParticles()**.

NOTE: Particle indices are 1-based, not 0-based!

# TIP:

For optimized particle access, use the "getAllXXX" functions to receive an array filled with all particle data for a particular property. Getting and iterating that array will be much faster than repeated calls to the property access functions for each particle by their index.

- [obj].getParticleID {index}: returns the unque ID for the specified particle index.
- [obj].getAllParticleIDs(): returns an array of all particle IDs for the prepared frame.
- [obj].getParticleAge {index}: returns the age (in frames) for the specified particle index.
- [obj].getAllParticleAges(): returns an array of all particle ages (in frames) for the prepared frame.
- [obj].getParticleTM {index}: returns the transform for the specified particle index.
- [obj].getAllParticleTMs(): returns an array of all particle transforms for the prepared frame.
- [obj].getParticlePosition {index}: returns the position for the specified particle index.
- [obj].getAllParticlePositions(): returns an array of all particle positions for the prepared frame.
- [obi].getParticleScale {index}: returns the scale for the specified particle index.

- [obj].getAllParticleScales(): returns an array of all particle scales for the prepared frame.
- **[obj].getParticleVelocity {index}**: returns the velocity (in units per frame) for the specified particle index.
- [obj].getAllParticleVelocities(): returns an array of all particle velocities for the prepared frame.
- [obj].getParticleShapeMesh {index}: returns the trimesh for the specified particle index.
- [obj].getAllParticleShapeMeshes(): returns an array of all particle trimeshes for the prepared frame.
- [obj].getParticleMatID {index}: returns the material ID override for the specified particle index.
- **[obj].getAllParticleMatIDs()**: returns an array of all particle material ID overrides for the prepared frame.

# **NOTE:** A material ID override value of 0 means no override is assigned to the particle.

- [obj].getParticleInstanceID {index}: returns the instance ID for the specified particle index.
- [obj].getAllParticleInstanceIDs(): returns an array of all particle instance IDs for the prepared frame.
- **[obj].getParticleSimGroups {index}**: returns the simulation group bitflags for the specified particle index.
- **[obj].getAllParticleSimGroups()**: returns an array of all particle simulation group bitflags IDs for the prepared frame.
- [obj].getParticleExportGroups {index}: returns the export group bitflags for the specified particle index.
- **[obj].getAllParticleExportGroups()**: returns an array of all particle export group bitflags IDs for the prepared frame.
- **[obj].getParticleUVWChannels {index}**: returns the mapping channel override indices for the specified particle index.
- **[obj].getAllParticleUVWChannels()**: returns an array of all particle mapping channel override indices for the prepared frame.
- **[obj].getParticleUVWs {index | channel}**: returns the mapping channel override value for the specified particle index.
- **[obj].getAllParticleUVWs {channel}**: returns an array of all particle mapping channel override values for the prepared frame.

# TIP:

Here is an example script, showing how to access individual particle transforms at frame 15, for a **tyFlow** object named "tyFlow001":

```
tf = $tyFlow001
tf.updateParticles 15
numParticles = tf.numParticles()
tms = tf.getAllParticleTMs()
for j in 1 to numParticles do
(
    tm = tms[j]
```

# tyFlow Modifiers

These are the tyFlow Modifiers, use the index or front quick link page to jump to each one.

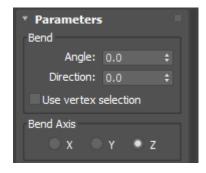


# tyBend Modifier

The **tyBend** modifier is an alternative to 3ds Max's built-in Bend modifier, that has options for bend segmentation so that you can bend a mesh along any axis and get a smooth/sharp crease that is not dependent on underlying topology.

# TIP:

The **tyBend** modifier is ideal for creating folding effects, like procedural origami.



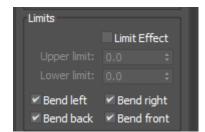
#### **Bend**

**Angle**: the amount (in degrees) to bend around the specified bend axis.

**Direction**: the amount (in degrees) to rotate the bend transform around the specified bend axis.

**Use vertex selection**: controls whether or not underlying vertex selections will affect the bend amount.

#### **Bend Axis**



**X/Y/Z**: the axis (of the bend gizmo) around which to bend geometry. Limits

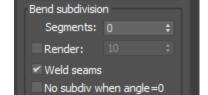
**Limit effect**: controls whether the bend deformation will be limited to a certain distance above/below the bend axis.

**Upper limit**: the amount to limit the bend deformation above the bend axis.

**Lower limit**: the amount to limit the bend deformation below the bend axis.

**Bend left/right/back/front**: these additional limits allow you to control whether bend deformations will happen on either sides of the two auxiliary axes of the bend transform.

**Bend subdivision** 



**Segments**: the number of segments to generate parallel to the slice plane.

**Render**: enables a render-only segment value, for increasing/decreasing segment count at render-time.

**Weld seams**: controls whether the vertices on the open edges of the segment slices will be welded together.

**No subdiv when angle=0**: when enabled, the mesh will only be subdivided when the bend angle is not equal to 0

# tyBoolean Modifier

The **tyBoolean** modifier is an experimental modifier capable of performing boolean CSG operations on a mesh. It can often provide results that contain less artifacts than boolean operations performed with 3ds Max's default boolean compound objects. It will also take the reference object's animated transformation changes into effect when performing its operations.



# **Input Objects (B)**

**Object list**: the list of objects that will act as object(s) B in the boolean operation.

NOTE: The object the modifier is applied to is object A.

**Hide after adding**: controls whether objects added to the list will be hidden after adding them.

# **Operation**

**Union**: joins object A and B together, into a single mesh.

**Intersection**: computes the intersection between object A and B as the resulting mesh.

**Subtraction** (**A - B**): subtracts object A from B and returns the resulting mesh.

**Subtraction** (**B** - **A**): subtracts object B from A and returns the resulting mesh.

**Slice** (**A - B**): subtracts object A from B (without capping subtracted faces) and returns the resulting mesh.

**Slice** (**B** - **A**): subtracts object B from A (without capping subtracted faces) and returns the resulting mesh.

# **Edge visibility**

**Coplanar tris -> polys**: controls whether groups of adjacent, coplanar triangles in the resulting mesh will be converted into polygons (have their shared edges hidden).

# **Pre-process**

**Split elements**: when enabled, the elements of the source mesh will be split apart and processed in parallel.

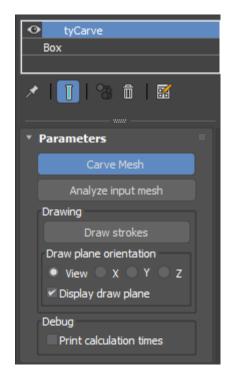
NOTE: If your mesh has shell surfaces (ie, a closed mesh with thickness), this option should be disabled. If this option is not disabled for shell surfaces, the inside and outside of the shell will be processed separately and may not return the desired result.

# **Post-process**

**Get orig element matID**: after the boolean operation is completed, the material ID from the first face of each original mesh element will be assigned to the output mesh element, as opposed to the material ID from the face of the operand mesh.

# tyCarve Modifier

The **tyCarve** modifier allows you to paint detailed slice patterns onto a mesh, which break up the mesh into concave chunks using CSG operations. The resulting chunks can be input into a **tyFlow** simulation as particles. When used to simulate destruction, these chunks have much more visual fidelity than the types of chunks created by a voronoi fracture operation.



#### **Parameters**

**Carve mesh**: enabling this checkbutton causes the modifier to evaluate. To disable modifier evaluation, uncheck this button.

**Analyze input mesh**: analyzes the input mesh (the mesh of the object the modifier is applied to), and returns information about issues in its topology.

NOTE: **tyCarve** works best on closed meshes that contain no self-intersections, coincident faces, degenerate faces, etc. Open meshes or meshes with topological problems can prevent **tyCarve** from performing clean cuts.

**Draw strokes**: enabling this checkbutton will enable stroke drawing in the viewport. Disabling this checkbutton will end the current stroke drawing session. In order to carve up a mesh, you must draw desired slice strokes over it.

**Plane orientation**: controls the orientation of the drawing plane. "View" means slices can be drawn in screen-space. "X/Y/Z" means slices will be drawn in world-space, along the chosen axis.

**Display draw plane**: controls whether the drawing plane is displayed as a visible grid in the viewport, during a stroke drawing session.

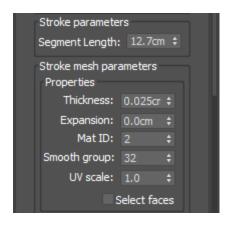
**Print calculation times**: controls whether details about the time taken to evaluate the carve are printed to the MAXScript Listener.

# **Strokes**

**Stroke list**: contains a list of drawn strokes that will be used to carve the mesh. A series of buttons below it can be used to enable/disable/remove/reorder strokes. When the modifier evaluates, strokes are evaluated in the top-to-bottom order that they appear in this list.

**NOTE: tyCarve** converts drawn strokes into 3D stroke meshes used to carve a mesh. The extrusion direction of a stroke is determined by the orientation of the plane it was drawn on.





**Segment length**: controls the size of segments generated between points on stroke meshes. Decreasing this value will increase the resolution of stroke meshes.

**Thickness**: controls the thickness of stroke meshes. Increase this value will create gaps between different elements of a carved meshes.

**Expansion**: controls the distance stroke meshes will be expanded outwards prior to carving. The direction of expansion is outward from the edges of the stroke plane.

**Mat ID**: controls the material ID of faces created to cap holes between carved elements.

**Smooth group**: controls the smoothing group value of faces created to cap holes between carved elements.

**UV scale**: controls the size of UVW coordinates generated on faces created to cap holes between carved elements.

**Select faces**: controls whether faces created to cap holes between carved elements will be selected in the resulting mesh.

**Displacement direction** 

**Explicit normals**: noise will displace stroke meshes along the direction of explicit normals on a stroke mesh.

**Average normals**: noise will displace stroke meshes along the direction of averaged normals on a stroke mesh.

**World Space**: noise will displace stroke meshes in world-space directions on a stroke mesh, determined by the resulting noise values.

Noise displacement 1 & 2

**Strength**: controls the strength of the noise displacement.

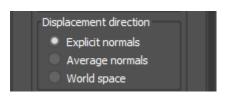
**Roughness**: controls the roughness of the noise displacement.

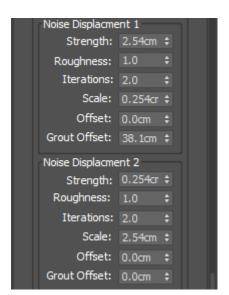
**Iterations**: controls the number of iterations the noise algorithm will undergo.

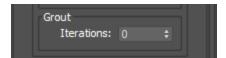
**Scale**: controls the scale of the noise displacement.

**Offset**: controls the offset assigned to input particle positions sent through the noise algorithm.

**Grout offset**: controls the progressive offset assigned to each grout iteration. The more grout iterations, the more offset.







#### Grout

**Iterations**: controls the number of grout iterations.

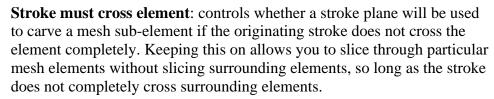
# NOTE:

"Grout" refers to the result of closely overlapping carve calculations, which can add more small carve details near stroke planes surfaces in a resulting mesh. For every grout iterations, all stroke planes will be reevaluated. Adding noise displacement to stroke meshes, and increasing grout offset will result in overlapping-but-not-coincident stroke planes being created to carve the mesh.

# **WARNING:**

For every grout iteration added, the entire stroke list must be reevaluated. Too many grout iterations can cause carve evaluation to slow down dramatically.

**Carving** 



**Bounds size** %: mesh elements whose bounding box size is less than this value (relative to the bounding box size of the whole mesh) will not be sliced. Increasing this value can reduce the number of slice operations performed on meshes with many smaller sub-elements, reducing total calculation time.

**Slice/subtract**: the "slice" operation will not cap the faces of sliced meshes, resulting in open mesh sub-elements. The "subtract" operation will cap the faces of sliced meshes, resulting in closed mesh sub-elements.

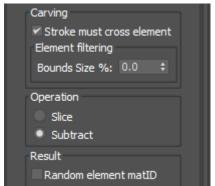
**Random element matID**: controls whether resulting mesh sub-elements will receive a randomized material ID. Turning this on and assigning a multi-sub material to the object with random colorings can help visualize the slices being drawn on a mesh.

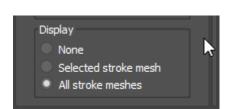
**Display** 

**None**: disables stroke mesh display in the viewport.

**Selected stroke mesh**: displays the stroke mesh selected in the stroke list in the viewport.

**All stroke meshes**: displays all stroke meshes in the viewport.





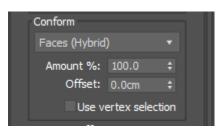
# tyConform Modifier

The **tyConform** modifier allows you to conform the surface of one piece of geometry to another.



#### **Parameters**

**Target Objects**: the list of objects to which the input surface will be conformed.



# Faces (Hybrid) Vertices Faces Faces (Hybrid) Expandwrap Shrinkwrap Raycast Matching Vertex ID Distance: 25-1000.000 v

#### Conform

**Vertices**: vertices of the input surface will conform to the nearest vertex in the target objects list.

**Faces**: vertices of the input surface will conform to the nearest face in the target objects list.

**Faces** (**Hybrid**): vertices of the input surface will conform to the nearest face in the target objects list (using a fast, hybrid algorithm based on nearest-vertex comparisons).

**Expandwrap**: vertices of the input surface will conform to the nearest hit point of a ray cast from the pivot of the nearest target object to each vertex.

**Shrinkwrap**: vertices of the input surface will conform to the nearest hit point of a ray cast from each vertex to the pivot of the nearest target object.

**Raycast**: vertices of the input surface will conform to the nearest hit point of a ray cast along a user-defined vector.

**Matching Vertex ID**: vertices of the input surface will conform to the matching vertex (by ID) of the nearest object in the target objects list.

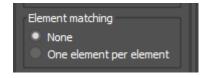
**Amount %**: the amount to conform the input surface to the target surface.

**Offset**: the offset from the target surface to apply the conform.

**Use vertex selection**: only selected vertices will be conformed (soft-selections are allowed).



**None**: disables element matching.



Distance Affect
Distance: 254000.0c \$
Falloff: 254.0cm \$

One element per element: when enabled, each element of the input surface will conform to only a single element of any mesh in the target objects list. This helps to reduce stretching artifacts, when vertices of a single input element conform to different elements of the target object list due to natural differences in their proximity. This setting is most useful when both inputs meshes and meshes in the target object list contain many different small elements.

**Distance Affect** 

**Distance**: vertices within this distance to the target surface will be fully affected.

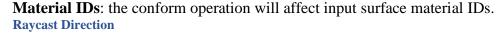
**Falloff**: the effect on vertices beyond the base distance to the target surface, but within this falloff distance, will diminish according to the inverse-square law.

Affect

**Vertices**: the conform operation will affect input surface vertex positions.

**Normals**: the conform operation will affect input surface vertex normal directions.

UVWs: the conform operation will affect input surface UVW coordinates.



**Vert to self pivot**: the raycast rays will be cast from the location of each vertex to the pivot of the source object.

**Absolute**: the raycast rays will be cast along the vector defined by the X/Y/Z values.

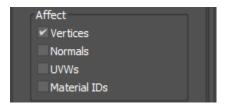
**Object z-axis**: the raycast rays will be cast along the local z-axis of a specified object.

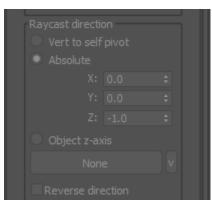
**Object**: the object whose z-axis will cast the rays.

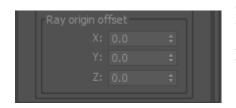
**X/Y/Z**: the coordinates of the user-defined vector to use as the raycast direction for Raycast mode.

**Reverse direction**: reverses the direction of the raycast rays. Ray Origin Offset

**X/Y/Z**: offset values to apply to each vertex's raycast origin point.

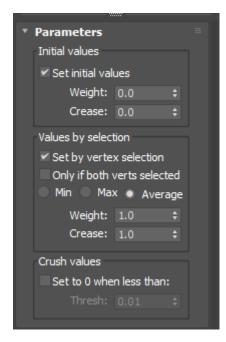






# tyEdgeWeights Modifier

The **tyEdgeWeights** modifier allows you to procedurally assign poly mesh edge weight/crease values (currently the only other way to assign these values is through manual entry in the Edit Poly modifier). These weights can be used to influence the results of other modifiers such as Meshsmooth, Quad Chamfer, etc.



# **Initial Values**

**Set initial values**: controls whether initial values will be assigned to all edges in the mesh.

Weight: the initial weight value.

**Crease**: the initial crease value.

# Values by selection

**Set by vertex selection**: controls whether values will be assigned to edges (after initial values are potentially set) based on the mesh's current vertex selection.

## INFO:

The assigned value will be the average selection value of each edge's two vertices. Vertex selection values are between 0 and 1, depending on soft selection, where 1 means the vertex is fully selected.

**Only if both verts selected**: edge weights will only be assigned to edges whose vertices are both selected.

**Min/Max/Average**: controls how the two selection values taken from each edge's vertices will be combined, to create the resulting edge weight/crease multiplier.

Weight: the initial weight value.

Crease: the initial crease value.

# **Crush values**

# INFO:

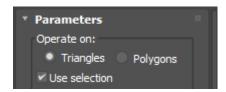
When using soft selections on a high resolution mesh, many edges can end up with tiny weight/crease values that will decrease the performance of certain modifiers that are dependent on edges with positive weight/crease values. Crushing small values to 0 will filter out those edges from other weight or crease-based algorithms, allowing them to run faster.

**Set to 0 when less than**: enables edge weight/crease value crushing, based on a threshold value.

**Thresh**: the crush threshold. If an edge's final weight value is below this threshold, its weight/crease values will be set to 0.

# tyFaceExtrude Modifier

The **tyFaceExtrude** modifier allows you to procedurally extrude and scale mesh faces, using vertex soft selection values to control the extrusion/scale amounts.

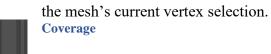


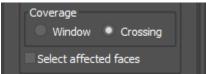
# Operate on

**Triangles/Polygons**: controls which mesh elements will be extruded.

NOTE: Polygons are groups of faces whose adjacent edges are invisible.

Use selection: controls whether extrusion/scale amounts will be based on





**Window**: all vertices of a face must be selected in order for the face to be extruded.

**Crossing** at least one vertex of the face must be selected in order for the face to be extruded.

**Select affected faces**: when enabled, affected faces (faces that are extruded or scaled) will be selected.

# **Face Extrude**

**Amount**: the amount to extrude applicable faces.

**Variation** %: the per-particle percentage of variation to apply.

**Along face normals**: extruded face vertices will be moved along the direction of the extruded face normal.

**Along face verts**: extruded face vertices will be moved along the direction of the extruded face vertex normals.

**Curve**: soft selection values will be interpolated along this curve.

**Relative to area**: the length of the extrusion will be relative to the area of the face.

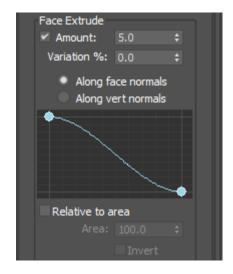
**Area**: the target face area.

**Invert**: inverts the effect of the face area multiplier.

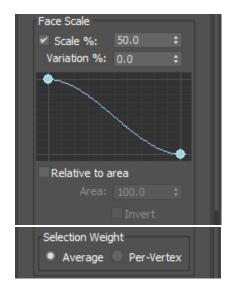
# **Selection Weight**

**Average**: the average selection weight of a face's vertices will be used to control the extrude amount.

**Per-Vertex**: the extrude amount of each face's vertex will be individually controlled by the vertex's selection weight.







# **Face Scale**

**Scale** %: the amount to scale applicable faces.

Variation %: the per-particle percentage of variation to apply.

**Curve**: soft selection values will be interpolated along this curve.

**Relative to area**: the amount to scale will be relative to the area of the face.

**Area**: the target face area.

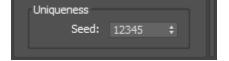
**Invert**: inverts the effect of the face area multiplier.

**Selection Weight** 

**Average**: the average selection weight of a face's vertices will be used to control the scale amount.

**Per-Vertex**: the scale amount of each face's vertex will be individually controlled by the vertex's selection weight.

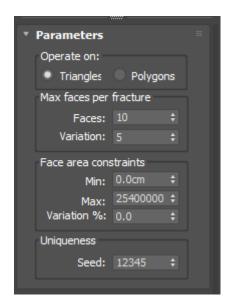
Uniqueness



**Seed**: the seed value for all varied parameters.

# tyFaceFracture Modifier

The **tyFaceFracture** modifier allows you to fracture mesh faces into separate mesh elements.



**NOTE:** The functionality of this modifier is equivalent to the functionality of **tyFlow's** Face Fracture operator.

# **Operate On**

**Triangles**: triangles will be treated as single faces.

**Polygons**: polygons (adjacent triangles that share a hidden edge) will be treated as single faces.

Max faces per fracture

**Faces**: the number of faces per fracture group.

**Variation**: the per-particle amount of variation to apply.

Face area constraints

Min/Max: the target min/max combined face area per fracture group.

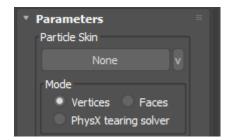
**Variation %:** the per-particle percentage of variation to apply.

Uniqueness

**Seed**: the seed value for all varied parameters.

# tyParticleSkin Modifier

The **tyParticleSkin** modifier can be used to deform a regular mesh with a **tyFlow** particle system.



# **Particle Skin**

**Node**: the **tyFlow** object whose particles will be used to skin the mesh.

#### Mode

**Vertices**: the mesh will search for nearest particles within the flow from the location of its vertices.

**Faces**: the mesh will search for nearest particles within the flow from the location of its face centers.

**PhysX tearing solver**: each individual vertex will be weighted to the nearest input particles. The PhysX-bind hierarchy of weighted particles will then be examined, and if breaks are detected over time, weightings will be adjusted to account for lost connections between affecting particles. The adjusted weightings manifest as gaps between previously-adjacent faces (ie, tears in the mesh). This effectively allows you to use **tyParticleSkin** as a history-independent tearing solver for rigid body simulations.

#### TIP:

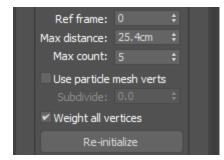
The **tyParticleSkin** modifier will perform the detach operation if it detects that the total number of mesh vertices is not equal to the total number of mesh faces multiplied by three (ie, the only valid ratio of faces-to-verts allowed for this mode), however, the modifier will perform faster if a **tyFaceFracture** modifier is placed below it, which does the proper detaching instead. In that configuration, the **tyFaceFracture** modifier (when operating on a static mesh) will only need to evaluate once, as opposed to the detach operation evaluating every frame when performed in the **tyParticleSkin** modifier.

#### NOTE:

Due to the requirement for detached faces by the PhysX tearing solver, when using that mode it is recommended to add a **tyWeld** modifier above the **tyParticleSkin** modifier in the modifier stack, so that detached faces are re-attached after the skinning operation is performed. This will have a performance cost but will keep resulting surfaces smooth instead of faceted.

# **INFO:**

If particles have a shape mesh, the vertices of the shape mesh will be considered during the nearest particle search, for improved accuracy.



**Ref frame**: the reference frame to use as the initialization time for the skin.

**Max distance**: the maximum distance that a particle can be from a mesh vertex/face in order to influence it.

**Max count**: the maximum number of particles that can influence a mesh vertex.

**Use particle mesh verts**: controls whether nearest-particle distances will be calculated to particle mesh vertices, instead of just particle positions.

**Subdivide**: controls whether randomly-placed implicit vertices will be added to the surface of applicable particle meshes, which can help to increase the accuracy of weight calculations.

Weight all vertices: if no particles are within the maximum distance to a mesh vertex/face, the nearest particle (regardless of distance) will be used. This ensures all vertices are weighted to at least one particle, assuming at least one particle exists.

# INFO:

During initialization, influencing particles will be weighted based on their relative and normalized distance to the target vertex. If a vertex is influenced by multiple particles, the weights for each particle will be determined by their distances to the vertex, with closer particles having a higher weight than further particles.

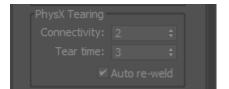
**Re-initialize**: re-initializes the skin.

# **Simulation Groups**

**Simulation groups**: controls which particle simulation groups will be imported.

# **Export Groups**

**Export groups**: controls which particle export groups will be imported.



Simulation Groups

Overlapping

Export Groups

1 2 3 4 5 6 78

9 10 11 12 13 14 1!16

Equal Not equal

ABCDEFCH

IJKLMNCP

Not equal

Overlapping

Equal

# **PhysX Tearing**

**Connectivity**: the maximum hierarchy depth to search for a PhysX bind connection between two particles which both influence a single vertex. Of a connection cannot be found within this depth, the particles are considered broken apart.

**Tear time**: the time, in frames, vertices will interpolate between bound and broken states after particle binding is found to be broken. The greater the value, the slower/smoother each new tear will form.

**Auto re-weld**: re-welds faces extracted by the **tyParticleSkin** modifier when in "PhysX tearing solver" mode. If this option is disabled, none of the mesh's individual faces will remain connected (the PhysX tearing solver requires faces to be separated prior to calculating tears).

# tyProperties Modifier

The **tyProperties** modifier allows you to assign **tyFlow**-specific properties to an object, which will be assigned to the corresponding particles derived from that object in a **tyFlow** simulation. For example, if you add an object to a Birth Objects operator that has a **tyProperties** modifier applied, the assignments enabled in the **tyProperties** modifier will transfer over to the particles created by that Birth Objects operator.



# **Simulation Groups**

**Set**: enables simulation group assignments.

**Simulation groups**: controls which particle simulation groups will be assigned.

# **Export Groups**

**Set**: enables export group assignments.

**Export groups**: controls which particle export groups will be assigned.

# **PhysX Binds**

**Bind setup ID**: controls whether a PhysX bind setup ID will be assigned.

#### **Custom Floats**

**Float list**: the list of custom floats which will be assigned.

Channel: the channel name of a custom float to assign.

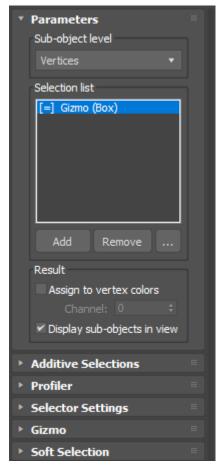
Value: the value of the custom float which will be assigned.

# tySelect Modifier

The **tySelect** modifier allows you to create procedural selections of mesh sub-objects (vertices, triangles, polygons or elements), using various customizable criteria.

**NOTE:** This modifier is meant to replace 3ds Max's default "Volume Select" modifier, which is out-dated and lacking a lot of important features.

# Parameters Rollout



**Sub-object level**: this controls what type of sub-object the **tySelect** modifier will select.

**Selection list**: this is the list of selectors that will be evaluated from top to bottom.

# INFO:

The **tySelect** modifier has its own selection list (similar to 3ds Max's modifier stack) which can be filled with selectors, each of which can have its own selection method and settings. The selection list is evaluated from top to bottom, and its purpose is to keep the **tySelect** modifier compact within 3ds Max's modifier stack. So if you need to create a selection with various criteria (adding and subtracting subobjects based on different selection methods), you'll only need 1 **tySelect** modifier in the modifier stack to do it, which keeps the modifier stack tidy in cases where complex selections are generated.

**Assign to vertex colors**: the resulting vertex selection values will be converted into color values and assigned to the mesh's vertex color map channel.

**Channel**: the channel to assign vertex colors to.

# TIP:

The default vertex color channel is 0, but technically any channel can take vertex color values.

**Display sub-objects in view**: controls whether the selected sub-object type will be displayed in the viewport (vertices, faces, etc).



#### INFO:

The additive selection parameters allow you to accumulate selections over time. The frames in the defined range will be evaluated and added together in succession, with their selections potentially decaying/fading as time progresses, depending on the temporal effect settings.

**Additive selections ENABLED/DISABLED**: this checkbutton controls whether or not additive selections will be read from the baked data.

# Frame Range

**Start/End**: the start and end frame of the additive selection interval.

**Steps/frame**: the number of time steps to sample/evaluate between frames during additive selection calculations.

**NOTE:** The more steps per frame, the more accurate the overall selection will be (temporally). Increasing the steps parameter is most useful for increasing accuracy when selections are generated from fast-moving objects.

## **Temporal Effect**

**Decay**: selection values will be decayed at this rate, over time.

**Age curve**: selection values will be interpolated along a curve by their age relative to a total lifespan value.

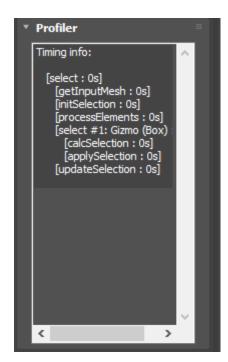
**Decay multiplier**: the multiplier applied to selection values over time.

**Lifespan**: the maximum lifespan of selection values.

**Bake additive selections**: computes the additive selections over the chosen frame range and stores them in memory.

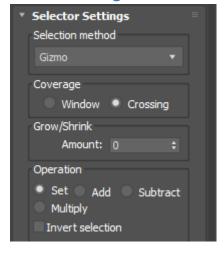
**Auto update**: additive selections will automatically be computed when changes to the input geometry or **tySelect** settings are detected.

# **Profiler Rollout**



The profiler rollout shows info related to the total calculation time for each step of the **tySelect** selection process.

# Selector Settings Rollout



**Selection Method**: controls the way in which mesh sub-objects will be selected for the selected selector in the selection list.

**Coverage (window/crossing)**: when window coverage is selected, all parts of a sub-object must be selected in order for the sub-object to be selected (for example, all vertices must be selected in order for a face to be selected, and all faces must be selected in order for an element to be selected). When crossing is selected, any parts of a sub-object can be selected in order for the sub-object to be selected.

**Grow/Shrink**: controls how must to grow or shrink a selection with adjacent sub-objects (grow with positive values, shrink with negative values).

# **Operation**

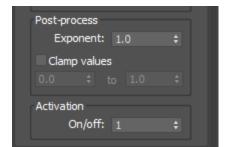
**Set**: sets the sub-object selection to the result of the selector.

**Add**: adds the result of the selector to the current selection.

**Subtract**: subtracts the result of the selector to the current selection.

**Multiply**: multiplies the result of the selector to the current selection.

**Invert**: inverts the result of the selector.



#### **Post-process**

**Exponent**: raises resulting soft-selection values to the specified power.

# INFO:

Changing the exponent setting from the default value of 1.0 is an easy way to crunch or inflate the influence amount of individual soft selection values.

**Clamp**: enables value clamping.

# **INFO:**

Certain operations can produce vertex selection values outside the range of [0-1]. Sometimes this can be intentional, but other times it can be an unintentional consequence of the selector's operation type. The clamp settings allow you to force resulting vertex selection values to be stay within a certain range.

**Min**: the minimum allowable vertex selection value after the selector's operation has completed.

**Max**: the maximum allowable vertex selection value after the selector's operation has completed.

#### Activation

**Activation**: an animatable parameter controlling the activation state of the selector (1 = on, 0 = off)

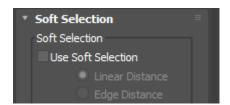
# TIP:

Set keyframes on the activation parameter to animate a selector's activation state over time.

# **Soft Selection Rollout**

# INFO:

Soft selections are only available in vertex sub-object mode, and emanate outwards from fully-selected vertices. Enabling soft selections will override any soft vertex selection values assigned directly by selectors.

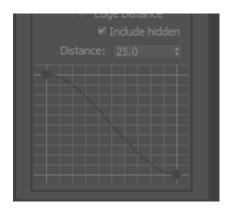


**Use Soft Selection**: enables soft-selection value calculations.

**Linear distance**: soft selection values will be computed as the distance from a vertex to the nearest fully-selected vertex, in worldspace.

**Edge distance**: soft selection values will be computed as the distance from a vertex to the nearest fully-selected vertex, along connected surface edges.

**NOTE:** When soft-selecting in "edge distance" mode, any vertices that have no direct connection along surface edges to fully-selected vertices will always receive a soft-selection value of 0.

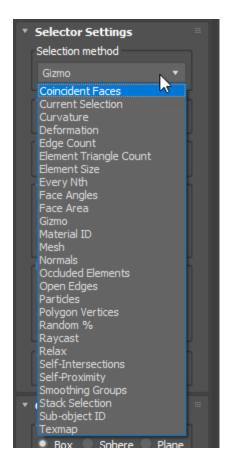


**Include hidden**: when enabled, the edge distance algorithm will traverse hidden edges. When disabled, it will skip them.

**Distance**: the distance value used to compute soft-selection ratios.

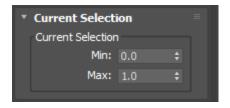
**Curve**: the curve used to interpolate soft-selection ratios along the distance, resulting in the final soft-selection value.

#### **Selection Methods**



Depending on the selection method, additional rollouts appear, they will be covered next few pages;

# **Current Selection**



The current selection selector imports the current selection stack as its selection. An example usage of this modifier, is if you want to re-compute soft selection values on a particular selection stack, in a way that is independent from the soft-selection methods previously computed. For example, the Mesh selector computes soft-selections by calculating surface proximities. This may not be ideal in situations where you want a soft-selection that is not relative to surface distances. So in that instance, you

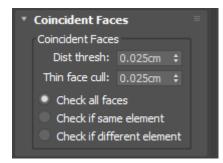
would use this selector to import the selection values of the Mesh selector, and re-compute the soft-selection.

Min: the minimum selection value to import.

Max: the maximum selection value to import.

NOTE: Values less than 1 are soft-selection values.

# **Coincident Faces**



The coincident faces selector computes selections by checking whether or not faces within a certain proximity to each other are overlapping.

**Dist thresh**: the maximum distance between two faces in order for them to perform an overlap test.

**Thin face cull**: faces with an angle between two edges smaller than this value will be ignored.

# TIP:

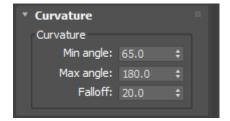
Very thin faces can return false-positive results in the overlap test. Culling thin faces like this will prevent them from being erroneously selected.

**Check all faces**: all faces will be compared to each other.

**Check if same element**: only faces of the same element will be compared to each other.

**Check if different element**: only faces of different elements will be compared to each other.

## Curvature



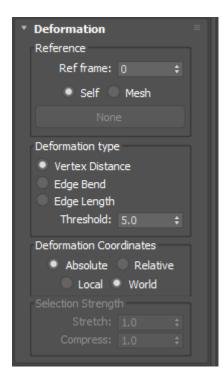
The curvature selector computes selections by measuring the angle between adjacent faces and comparing the result to user-defined threshold parameters.

**Min/Max angle**: faces whose adjacency angle is between these values will be fully selected.

**Falloff**: faces whose adjacency angle is outside of the min/max threshold but within this falloff threshold will have their vertices partially selected, depending on how close their angle is to the min/max threshold.

#### **Deformation**

The deformation selector computes selections by comparing the input mesh to a reference mesh at a certain point in time. The greater the distance between mesh vertices, the stronger the selection. For example, it can be used to compute selections for stretched or creased areas of an animated mesh.



#### Reference

**Ref frame**: the frame at which the reference mesh will be analyzed.

**Self**: the reference mesh will be an internal snapshot of the current mesh evaluated at the reference frame.

**Mesh**: the reference mesh will be the mesh of an external object evaluated at the reference frame.

**NOTE:** Deformations can only be computed on meshes with identical topology. If the meshes do not have identical vertex counts, the deformation selection cannot be performed.

# **Deformation type**

**Vertex Distance**: the selection will be computed by comparing distances between vertices of the reference/input meshes.

**Edge Bend**: the selection will be computed by comparing angles between edges of the reference/input meshes.

**Edge Length**: the selection will be computed by comparing differences between the lengths of edges in the reference/input meshes.

**Threshold**: the type-dependent selection threshold. For example, it is used as an angle threshold in "edge bend" mode, and a length threshold in "edge length" mode.

## **Deformation coordinates**

**Absolute/relative**: controls whether the measured differences will be absolute values in the selected spatial coordinates, or values relative to adjacent vertex neighbors.

**Local/space**: controls whether differences will be measured in local (mesh) space coordinates, or world space.

## **Selection strength**

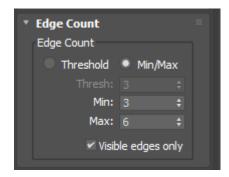
Selection strength values are available in "edge length" mode.

**Stretch**: a multiplier which controls the amount that stretched edges will contribute to the selection.

**Compress**: a multiplier which controls the amount that compressed edges will contribute to the selection.

# **Edge Count**

The edge count selector selects sub-objects based on how many adjacent edges they are connected to.

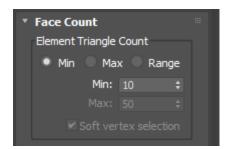


**Threshold**: a soft-selection will be generated based on the number of adjacent edges compared to a threshold value.

**Min/Max**: elements with an adjacent-edge count within the min/max range will be selected.

# **Element Triangle Count**

The element triangle count selector selects sub-objects based on how many triangles are in the sub-object's mesh element. A mesh element is a contiguous set of triangles.



**Min**: counts equal to the min value and above will be fully selected.

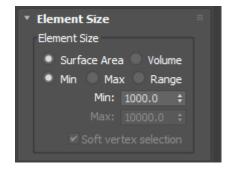
**Max**: counts equal to the max value and below will be fully selected.

**Range**: counts equal to the min value and above, or equal to the max value and below will be fully selected.

**Soft vertex selection**: elements with counts that didn't qualify for a full selection will be soft-selected based on the ratio between their count and the count criteria.

#### **Element Size**

The element size selector selects sub-objects based on the size of their mesh element. A mesh element is a contiguous set of triangles.



The element size selector selects sub-objects based on the size of their mesh element. A mesh element is a contiguous set of triangles.

**Surface Area/Volume**: the element size type.

**Min**: sizes equal to the min value and above will be fully selected.

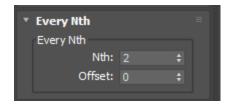
**Max**: sizes equal to the max value and below will be fully selected.

**Range**: sizes equal to the min value and above, or equal to the max value and below will be fully selected.

**Soft vertex selection**: elements with sizes that didn't qualify for a full selection will be soft-selected based on the ratio between their size and the size criteria.

# **Every Nth**

The every nth selector selects sub-objects based on repeating patterns. For example, an nth value of "2" will select every other sub-object. An nth value of "3" will select every third sub-object, etc.



**Nth**: the pattern value to use.

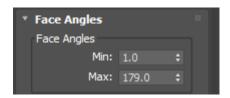
**Offset**: the offset to apply to the input sub-object index, prior to pattern matching.

# INFO:

The every nth selector uses the modulus operator to compute pattern matches. The selection condition equation is: (([sub-object index] + offset) % nth) == 0

# **Face Angles**

The face angles selector computes selections based on the angles between the edges of mesh faces. If all three pairs of adjacent edges within the same face have an angle within the defined min/max thresholds, the face will be selected.

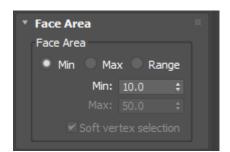


**Min**: the minimum angle between face edges in order for a face to be a candidate for selection.

**Max**: the maximum angle between face edges in order for a face to be a candidate for selection.

## Face Area/Volume

The face area selector selects sub-objects based on the area of mesh faces.



**Min**: areas equal to the min value and above will be fully selected.

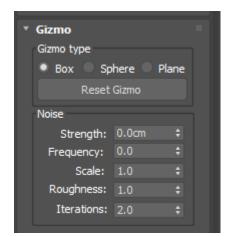
**Max**: areas equal to the max value and below will be fully selected.

**Range**: areas equal to the min value and above, or equal to the max value and below will be fully selected.

**Soft vertex selection**: faces with areas that didn't qualify for a full selection will be soft-selected based on the ratio between their area and the area criteria.

#### Gizmo

The gizmo selector selects sub-objects using a gizmo sub-object. Sub-objects encompassed by the gizmo sub-object will be fully selected.



Box/Sphere/Plane: the shape of the gizmo sub-object.

# Noise

The noise settings allow you to offset the way in which vertices are selected by the gizmo sub-object.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will increase selection evolution over time.

Scale: the scale of the noise.

**Roughness**: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

# **Material ID**

The material ID selector selects sub-objects using face material ID values.



**IDs**: the list of material IDs to match.

#### Mesh

The mesh selector selects sub-objects using the meshes of external scene objects.

## TIP:

Enable "use soft selection" in the soft selection rollout in order to interpolate mesh proximity values as vertex soft-selection values.



**Select if inside**: vertices inside the input meshes will be fully selected.

**Accuracy**: the accuracy of the inside test. Increase this value for meshes that are self-intersecting or have holes in them.

#### **Noise**

The noise settings allow you to offset the way in which vertex positions are measured during the distance/volume tests.

**Strength**: the strength of the noise.

**Frequency**: the frequency of the noise. Increasing this value will evolve the noise pattern over time.

Scale: the scale of the noise.

**Roughness**: the roughness of the noise.

**Iterations**: the number of iterations to use to calculate the noise.

# **Normals**

The normals selector selects sub-objects based on the direction of mesh face normals.



**Gizmo direction**: the direction vector of the gizmo sub-object determines which face normals will be selected.

**Absolute X/Y/Z**: the vector defined by the X/Y/Z spinners will determine which face normals will be selected.

**Towards/Away/Both**: controls the manner in which face normals must align to the test vector in order to be selected.

**Threshold**: faces with normals which align to the test vector within this threshold will be selected. The greater the threshold value, the less exact an alignment must be in order for a face to be considered a selection candidate.

#### **Occluded Elements**

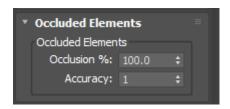
The occluded elements selector selects sub-objects based on whether or not all vertices of a mesh element are occluded by other parts of the mesh.

# TIP:

This selector makes it easy to select mesh elements that are inside of other mesh elements, or mesh elements whose normals are flipped inside-out. An example usage would be selecting all interior faces/elements of a blobmesh created by a **tvMesher**.

# INFO:

The occlusion test is performed by casting rays out from every vertex of the mesh. If a vertex doesn't hit a face, it is deemed un-occluded by the mesh. Increasing the accuracy parameter increases the number of rays cast from each vertex, which increases the probability that a vertex which is not completely occluded by parts of the mesh will have rays which escape without hitting a face.



**Occlusion** %: the minimum percent of vertices within a mesh element that must be occluded by parts of the overall mesh, in order for the element to be selected.

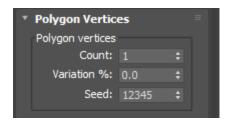
**Accuracy**: the number of rays to cast from each vertex, to test for occlusion.

# **Open Edges**

The open edges selector selects sub-objects based on whether or not they are touching an open edge. An open edge is an edge with only one adjacent face. This selector has no settings.

# **Polygon Vertices**

The polygon vertices selector selects a certain number of each polygon's vertices.



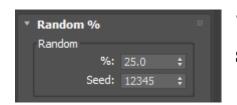
**Count**: the number of vertices in each polygon to select.

**Variation** %: the per-particle percentage of variation to apply.

**Seed**: the seed value which controls the variation.

# Random %

The random % selector selects a random percentage of sub-objects.



%: the random percentage of sub-objects to select.

**Seed**: the seed value which controls the randomness.

#### **Particles**

The particles selector selects sub-objects using the particles of external particle systems



The particles selector selects sub-objects using the particles of external particle systems.

**Absolute radius**: sub-objects within an absolute radius to particles will be selected.

**Radius**: the absolute radius value to use.

**Particle radius**: sub-objects within each particle's radius will be selected.

Multiplier: a multiplier to apply to each particle's radius.

**Simulation Groups** 

**Simulation groups**: controls which particle simulation groups will be imported.

**Export Groups** 

**Export groups**: controls which particle export groups will be imported.

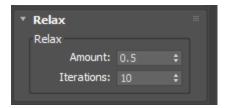
# **Raycast**

The raycast selector selects sub-objects based on whether or not faces are hit with rays emanating from a gizmo sub-object.



#### Relax

The relax selector applies a relaxation kernel to current vertex selections, which averages selection values between edge-adjacent vertices.



**Amount**: the amount to linearly-interpolate current selections to post-relax selection values.

**Iterations**: the number of times to run the relaxation kernel over the current selection.

## **Self-Proximity**

The self-proximity selector computes selections based on the distance of vertices to other vertices within the same mesh.



**Distance**: vertices within this distance to each other will be fully selected.

**Falloff**: vertices whose distance is further than the distance threshold, but within the falloff threshold, will be partially selected.

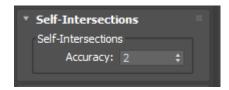
**Ignore same element**: vertices within the same mesh element will ignore each other.

**Ignore adjacent vertices**: vertices whose edge connectivity is within a certain depth threshold to each other will ignore each other.

**Depth**: the vertex connectivity depth threshold.

#### **Self-Intersections**

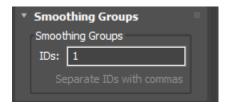
The self-intersections selector computes selections based on whether or not vertices are found to be intersecting their mesh.



**Accuracy**: controls the accuracy of the raycaster used to compute information about whether a vertex is inside or outside of its mesh.

### **Smoothing Groups**

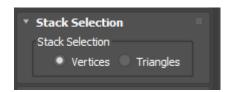
The smoothing groups selector selects sub-objects using face smoothing group values.



**IDs**: the list of smoothing group to match.

### **Stack Selection**

The stack selection selector selects sub-objects using existing selections present in the modifier stack.

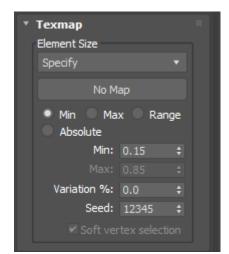


**Vertices**: existing vertex selections will be used to select sub-objects.

**Triangles**: existing triangle selections will be used to select sub-objects.

### **Texmap**

The texmap selector selects sub-objects using the monochromatic intensity values of a texmap.



**Texmap**: the texmap to sample per-vertex values from.

**Min**: texmap values equal to the min value and above will be fully selected.

**Max**: texmap values equal to the max value and below will be fully selected.

**Range**: texmap values equal to the min value and above, or equal to the max value and below will be fully selected.

**Absolute**: texmap values will be converted directly into selection values.

**Variation** %: the per-particle percentage of variation to apply.

**Seed**: the seed value which controls the variation.

**Soft vertex selection**: vertices with texmap values that didn't qualify for a full selection will be soft-selected based on the ratio between their texmap value and the texmap value criteria.

# tySlice Modifier

The **tySlice** modifier is an optimized alternative to 3ds Max's built-in Slice modifier, that has options for capping slice holes, multi-object slice plane input, etc.

## tySlice Rollout



#### **Slice Plane**

**Gizmo**: the slice will be controlled by the modifier's sub-object gizmo.

**Object**: the slices will be controlled by the z-axis of input object transforms.

**Object list**: the list of input objects to use as slice planes.

## **Slice operation**

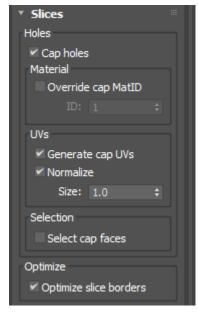
**Split**: the input mesh will be split in half along the slice plane(s).

**Refine**: the input mesh will be split along each refinement slice plane.

**Remove top**: the input mesh will be split along the slice plane(s) and any vertices above the slice plane will be removed.

**Remove bottom**: the input mesh will be split along the slice plane(s) and any vertices below the slice plane will be removed.

## Slices Rollout



#### Holes

**Cap holes:** controls whether slices will be capped with new faces.

### Material

**Override cap MatID**: controls whether cap faces will be given a material ID override.

**ID**: the cap face material ID value.

#### IIV

**Generate cap UVs**: controls whether UVW coordinates will be generated on new cap faces.

**Normalize**: controls whether cap UVW coordinates will be normalized.

**Size**: the size of the cap face UVW coordinates.

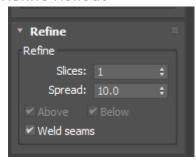
#### **Selection**

**Select cap faces**: controls whether new cap faces will be flagged for selection.

## **Optimize**

**Optimize slice borders**: controls whether slice borders will be optimized, removing things like extraneous vertices along parallel edges.

## **Refine Rollout**



**Slices**: the number of refinement slices.

**Spread**: the distance between refinement slices.

**Above/Below**: controls whether refinement slices will be created above/below the original slice plane(s).

**Weld seams**: controls whether the edges between refinement slices will be welded.

# tySmooth Modifier

The **tySmooth** modifier is an optimized and multi-threaded version of 3ds Max's built-in smooth modifier.



### Smooth

**Auto Smooth**: When disabled, no mesh faces will share smoothing groups. When enabled, mesh faces will be assigned smoothing groups based on the angles between their adjacent faces.

**Prevent indirect smoothing**: re-processes smoothing groups generated by the auto-smooth function, in order to prevent indirect smoothing.

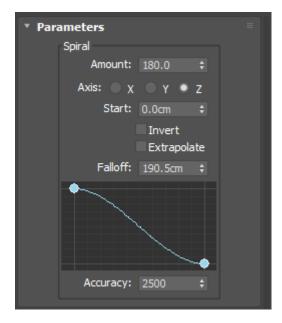
### **INFO**:

Indirect smoothing happens when adjacent faces (which do not satisfy the angle threshold condition) end up as members of the same smoothing group because they are connected by other adjacent faces which do satisfy the angle threshold condition. "Prevent indirect smoothing" attempts to rectify this problem by shuffling smoothing groups in a way that keeps properly-smoothed adjacent faces as members of the same group, while keeping indirectly-smoothed adjacent faces as members of different groups. Enabling this feature has a performance cost that will cause the auto-smooth algorithm to run slower.

**Threshold**: the angle threshold to use, to compare adjacent faces. If the angle between adjacent faces is less than this value, they will be assigned to the same smoothing group.

# tySpiral Modifier

The **tySpiral** modifier applies a circular deformation to geometry, outward from a central axis. It is similar to 3ds Max's built-in Twist modifier, except that instead of deforming geometry around an axis in an absolute manner, the strength of the circular effect has a falloff, allowing the deformation to form a spiral pattern.



### **Spiral**

**Amount**: the amount, in degrees, of spiral deformation to apply.

**Axis X/Y/Z**: the local axis of the sub-object gizmo around which to apply the spiral deformation.

**Start**: the minimum distance from the specified axis a vertex must be in order to undergo deformation.

**Invert**: inverts the spiral effect, so that vertices within the "start" distance will undergo deformation, rather than vertices further than the "start" distance.

**Extrapolate**: extrapolates the spiral deformation effect amount, for vertices that are further from the "start" distance to the specified axis than the falloff amount.

**Falloff**: the maximum distance past the "start" distance from the specified axis to apply the deformation.

**Falloff curve**: defines the deformation amount from "start" distance to ["start" + falloff] distance from the specified axis.

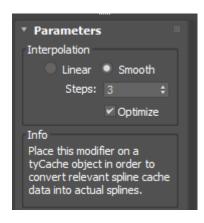
**Accuracy**: defines the internal resolution of the falloff curve. Values that are too low can produce geometry artifacts when the spiral amount value is high.

# tySplineCache Modifier

The **tySplineCache** modifier extracts spline cache data loaded into a **tyCache** object, and converts it into editable splines.

### INFO:

Spline data exported from an Export Particles operator is stored in a mesh-based format. This modifier, when placed on a **tyCache** object that contains that data, will convert it into a format that 3ds Max can recognize as traditional, editable splines.



## Interpolation

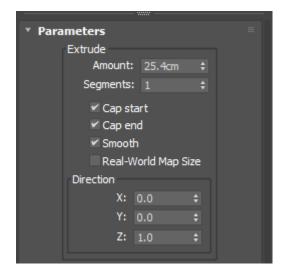
**Linear/Smooth**: the type of interpolation to apply to spline knots.

**Steps**: the number of interpolation steps to apply to splines.

**Optimize**: controls whether or not the step count will be optimized depending on the spline curvature.

# tySplineExtrude Modifier

The **tySplineExtrude** modifier is a simplified version of 3ds Max's built-in extrude modifier, with the key difference being that **tySplineExtrude** will properly apply underlying spline material IDs to the each whole extrusion (sides and cap), rather than just the *sides* of an extrusion.



### Extrude

**Amount**: the height of the extrusion.

**Segments**: the number of height segments that will span the extrusion.

**Cap start**: controls whether the start of the extrusion will be capped.

**Cap end**: controls whether the end of the extrusion will be capped.

**Smooth**: controls whether the side of the extrusion will be smoothed.

**Real-World map size**: controls whether UVW coordinates applied to the resulting mesh will be given a real-world size.

### **Direction**

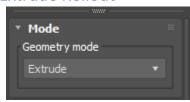
**X/Y/Z**: used to define the direction vector of the extrude operation.

# tySplineMesher Modifier

The **tySplineMesher** modifier can quickly convert huge numbers of splines into geometry, much faster than 3ds Max's default renderable spline modifier. It is ideal for **tySplines** objects, but can be assigned to shape objects of any type.

**NOTE:** Variation parameters are seeded by both the seed spinner values, and the material ID of the subspline. In order to apply variation between subsplines, ensure the subsplines have different material ID values.

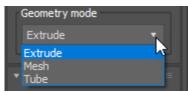
### **Extrude Rollout**



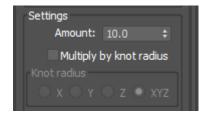
**Extrude**: extrudes the splines in a specific direction.

**Mesh**: deforms meshes along the length of splines.

**Tube**: converts splines into cylindrical meshes.



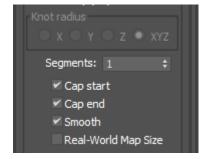
**Settings** 



**Amount**: the distance to extrude the splines.

**Multiply by knot radius**: controls whether the distance of spline extrusions is affected by knot radius. Knot radius is a custom value passed up the modifier stack by **tySplines** objects, based on originating particle radii. This option will not work if this modifier is assigned to a default 3ds Max shape.

**Knot Radius** 



**X/Y/Z**: the radius of the knot along the given axis will be the multiplier.

**XYZ**: the radius of the longest knot axis will be the multiplier.

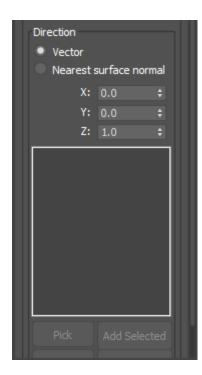
**Segments**: the number of segments to generate along the extrusion.

**Cap start**: faces will be generated on the bottom of closed extruded splines.

**Cap end**: faces will be generated at the top of closed extruded splines.

**Smooth**: extruded faces will be assigned smoothing groups.

**Real-world map size**: assigned UVW coordinates will be relative to the size of the extruded mesh.



**Vector**: the direction of the extrusion will be a user-specified vector.

**Nearest surface normal**: the direction of the extrusion will be the nearest surface normal on the nearest mesh to each input spline knot.

**Objects**: the list of input objects whose distance to spline knots will be measured.

### NOTE:

Capping options are not available when "nearest surface normal" direction mode is activated.

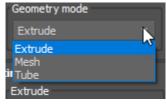
## Spline Mesher Rollout



### **Geometry mode**

**Tube**: converts splines into cylindrical meshes.

**Mesh**: deforms meshes along the length of splines.



### **Basic Settings**

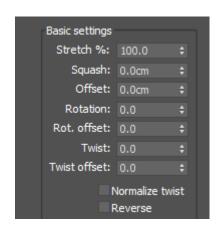
**Stretch** %: controls the maximum percent along splines to stretch geometry. Animating this value allows you to animate the length of an extrusion along the spline.

**Squash**: squashes the extrusion along a vector that is perpendicular to each spline segment.

**Offset**: offsets the extrusion along a vector that is perpendicular to each spline segment.

**Rotation**: rotates the extrusion along each spline segment's local axis.

**Rotation offset**: controls the starting point along the spline after which rotation values will take effect.



**Twist**: controls the amount of twist that will be applied to successive geometry cross sections along the length of the spline.

**Twist offset**: controls the starting point along the spline after which twist values will take effect.

**Normalize twist**: controls whether the amount of twist applied is dependent on the length of the spline.

**Reverse**: controls whether the ordering of knots along a spline is reversed.



### **Orientation**

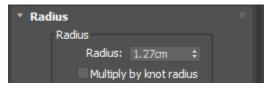
**From spline**: the orientation of geometry cross sections along the spline will be dynamically determined by the spline's topology.

**From spline**: the orientation of geometry cross sections along the spline will be static.

**From knots**: the orientation of geometry cross sections along the spline will be determined by its knots. This option will not work if this modifier is assigned to a default 3ds max shape.

**X/Y/Z**: controls which axis of the source TM will be the forward vector of the orientation matrix.

## Radius Rollout



**Radius**: controls the overall radius of spline extrusions.

**Multiply by knot radius**: controls whether the radius of spline extrusions is affected by knot radius. Knot radius is a custom value passed up the modifier stack by **tySplines** objects, based on originating particle radii. This option will not work if this modifier is assigned to a default 3ds max shape.



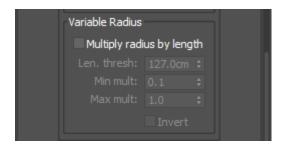
### **Knot Radius**

**X/Y/Z**: the radius of the knot along the given axis will be the multiplier.

**XYZ**: the radius of the longest knot axis will be the multiplier.

### Variable radius

**Multiply radius by length**: controls whether the radius of spline extrusions will be affected by the overall length of each spline.



**Length threshold**: controls the reference length that each spline length will be compared against, when determining how to multiply its extrusion radii.

**Min mult**: the minimum multiplier value for extrusion radii. Increase this value to prevent variable-radius spline extrusions from becoming too thin.

**Max mult**: the maximum multiplier value for extrusion radii. Decrease this value to prevent variable-radius spline extrusions from becoming too thick.

**Invert**: inverts the variable-radius operation.

## **Taper**

**By Length**: the length of the spline will determine the taper interpolation.

By Age: the age of the knots will determine the taper interpolation.

**Linear**: linear taper is a radius multiplier applied evenly along the spline.

**Curve**: curve taper is a radius multiplier applied by projecting a curve along the spline.

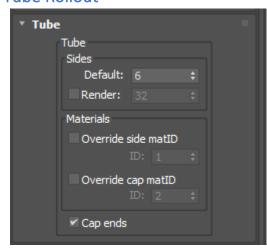
**Amount**: the amount of linear taper to apply.

**X/Y/XY**: controls the taper axis.

**Symmetrical taper**: enabling this option tapers splines from the center outwards, instead of from end-to-end.



## **Tube Rollout**



### **Sides**

**Default**: the default number of side faces created around a spline extrusion. Increasing this value increases the resulting mesh resolution.

**Render**: enabling this override allows you to control the number of side faces created around a spline extrusion at render time.

### **Materials**

**Override side matID**: controls whether side faces will receive a custom material ID.

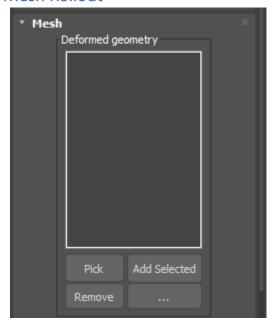
**Side ID**: controls the material ID value for side faces.

**Override cap matID**: controls whether cap faces will receive a custom material ID.

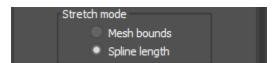
Cap ID: controls the material ID value for cap faces.

**Cap ends**: controls whether the ends of an extrusion are capped with faces.

## Mesh Rollout



Mesh list: the list of input meshes that will be randomly selected and extruded along input splines.



#### Stretch mode

Mesh bounds: the "stretch" parameter in the Spline Mesher rollout will act as a multiplier on the bounds of the input mesh. A value of 100 will stretch the input mesh to its default length.

Spline length: the "stretch" parameter in the Spline Mesher rollout will act as a multiplier on the length of the input spline. A value of 100 will stretch the input mesh to the length of the spline.

## Offset along spline

Applies an offset to the input mesh's position along the length of the spline.

**Percent**: the mesh will be offset by a percentage of the spline's overall length.

**Distance**: the mesh will be offset by an absolute value along the spline's overall length.

**Variation** %: the per-particle percentage of variation to apply.



Axis

**X/Y/Z**: the axis used to orient the input mesh along the spline.

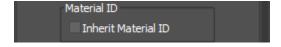
**Flip**: flips the orientation axis.

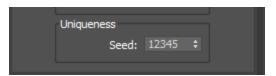
**Center pivot**: input meshes will be deformed along their center, rather than their default pivot. **Material ID** 

**Inherit Material ID**: input meshes will inherit the material ID of their spline.





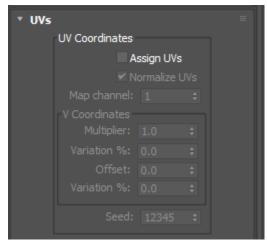




### Uniqueness

**Seed**: the seed value for all varied parameters.

## **UV Coordinates Rollout**



**Assign UVs**: (mesh mode only) controls whether UV overrides will be applied.

**Normalize UVs**: controls whether UVW coordinates assigned to a vertex are relative to that vertex's location along the spline.

### INFO:

If "normalize UVs" is enabled, texture vertex V-coordinates will have a value between 0.0 and 1.0, depending on how far along each spline they are. If "normalize UVs" is disabled, texture vertex V-coordinates will have a value relative to the number of edge segments along the subspline they are. For example, in "Tube" mode a vertex placed along the spline at the 6th edge segment will have a V-coordinate value of 6.0.

**Map channel**: (mesh mode only) controls the mapping channel that UV overrides will be applied to.

**NOTE:** U Coordinate settings are only available in "Tube" mode.

**Multiplier**: a multiplier applied to U coordinate values.

**Variation** %: the per-particle percentage of variation to apply.

**Offset**: an overall offset applied to U coordinate values.

**Variation** %: the per-particle percentage of variation to apply.

## V Coordinates

**Multiplier**: a multiplier applied to V coordinate values.

**Variation** %: the per-particle percentage of variation to apply.

**Offset**: an overall offset applied to V coordinate values.

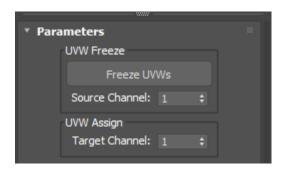
**Variation** %: the per-particle percentage of variation to apply.

**Seed**: the value which will seed the various randomization functions

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# tyUVWFreeze Modifier

The **tyUVWFreeze** modifier can be used to freeze existing UVW coordinates applied to an object at a particular time and map channel. It functions similar to how a default 3ds Max UVWUnwrap modifier will lock UVW coordinates in place, once applied, however the **tyUVWFreeze** modifier operates more efficiently, and frozen coordinate data is not lost if the underlying mesh topology changes.



#### **UVW Freeze**

**Freeze UVWs**: when this button is pressed, the UVW coordinates for a particular source channel will be frozen at the current frame.

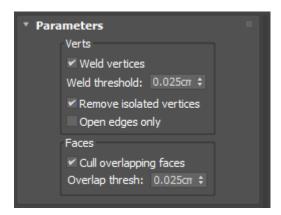
**Source channel**: the source map channel to retrieve the UVW coordinates from, on the input mesh.

## **UVW** Assign

**Target channel**: the target map channel where the frozen UVW coordinates will be assigned.

# tyWeld Modifier

The **tyWeld** modifier can be used to weld vertices and cull overlapping faces of a mesh. It is especially helpful for welding/culling the interior faces of a mesh converted to tets with a **tyMesher**.



#### Verts

**Weld vertices**: controls whether vertices will be welded together by proximity.

**Weld threshold**: the maximum distance threshold between vertices in order for them to be welded together.

**Remove isolated vertices**: removes any vertices not connected to at least one face.

**Open edges only**: only vertices on open edges (edges connected to a single face) will be welded.

#### **Faces**

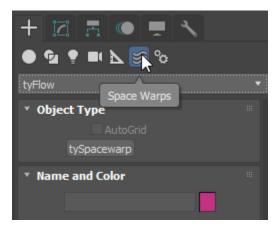
**Cull overlapping faces**: controls whether overlapping faces (faces whose verts are all coincident) will be removed from the mesh.

**Overlap thresh**: the maximum distance between vertices on two candidate faces in order for them to be considered coincident. If all three vertices are coincident, the faces are considered overlapping.

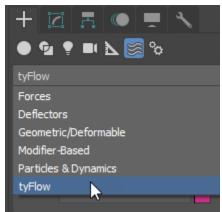
# **Spacewarps**

# tySpacewarp Object

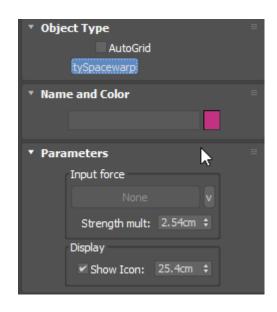
The **tySpacewarp** object can be used to convert **tyFlow** force helpers (ex: **tyVortex**) into spacewarps that are compatible with 3rd party plugins (ex: PhoenixFD).



Accessed from the Space Warps Tab



Select tyFlow from the Space Warps Menu



### **Input force**

**Force object**: the input **tyFlow** force helper which will be converted into a regular spacewarp.

**Strength mult**: a multiplier which will be applied to the resulting force calculations.

## **Display**

**Show icon**: controls whether or not to display the **tySpacewarp** icon in the viewport.

**Icon size**: controls the size of the viewport icon.

# **Splines**

# tySplines Object

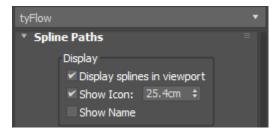
The **tySplines** object takes input particles sent from a **Spline Paths** operator and converts them into native 3ds max splines.

### NOTE:

You must select a **tySplines** object within a **Spline Paths** operator inside of an active flow in order to feed a **tySplines** object its data. A **tySplines** object contains a cache of data for fast playback, but does not save that cache to disk. Thus, its parent flow must always exist in the scene with it, in order for it to work.

### **WARNING:**

Multiple operators within the same flow can send particles to the same **tySplines** object, but you should not send particles from multiple flow objects into the same **tySplines** object, or else the resulting splines will not display correctly.



**Spline Paths** 

### **Display**

**Display splines in viewport**: controls whether or not to generate splines in the viewport.

**Show icon**: controls whether or not to display the **tySplines** icon in the viewport.

**Icon size**: controls the size of the viewport icon.

**Show name**: displays the name of the **tySplines** object next to its icon in the viewport.



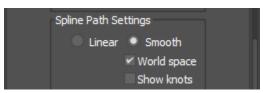
**Linear/Smooth**: controls the interpolation method used to generate spline.

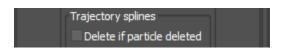
**World space**: controls the coordinate system to generate splines in. Turning this off will generate splines whose position is relative to the position of the **tySplines** object in the world.

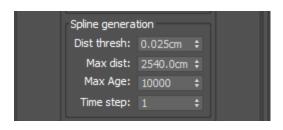
**Show knots**: controls whether or not to display each spline's knots in the viewport.

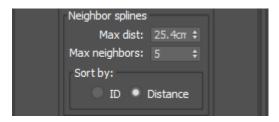
**Trajectory splines** 

**Delete if particle deleted**: trajectory splines whose originating particle is deleted will not generate the leftover trail that is within the trajectory/sibling max age limit.

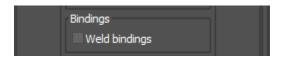


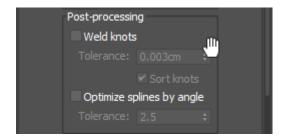












### **Spline generation**

**Dist thresh**: controls the minimum step size between points on a particle's trajectory in order for them to form a spline segment.

**Max dist**: controls the maximum distance between two sibling particles in order for them to be conjoined with a spline segment.

**Max age**: controls the maximum age of particle trajectory points in order for them to be adding to a spline segment.

**Time step**: controls the minimum time step between particle positions in order for them to form a spline segment.

## **Neighbor splines**

**Max dist**: controls the maximum distance between two particles in order for them to be conjoined with a spline segment.

**Max neighbors**: controls the maximum number of spline segments a particular particle can have, connecting it to its neighbors.

**Sort by ID/distance**: controls the method used to sort neighbors within the distance threshold for any particular particle, if the number of neighbors is larger than the value of "Max neighbors".

**Neighbor Simulation groups**: controls which particle simulation groups will be considered for the neighbor collection. **Bindings** 

**Weld bindings**: controls whether or not to weld spline segments corresponding to connected bindings between particles together, end-to-end.

### **Post Processing**

Weld knots: controls whether or not to weld resulting splines segments together, based on whether or not their endpoints are coincident.

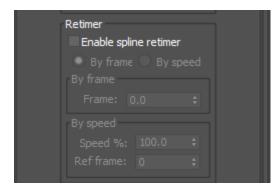
**Tolerance**: the weld distance tolerance. Spline endpoints whose distance is below this value will be welded together.

**Sort knots**: sorts the candidate knots prior to welding.

**NOTE:** Knot sorting has a performance cost, but keeping knots in order can help keep weld order deterministic between frames where knot count changes.

**Optimize splines by angle**: controls whether spline knot counts will be reduced based on relative angles between spline segments.

**Tolerance**: the tolerance angle for performing knot reductions. Spline segments whose relative angle is less than this value will have their adjacent knot removed.



#### Retimer

**Enable spline retimer**: controls whether or not to enable spline animation retiming.

**Retime type**: controls whether the retimer affects playback frame or speed.

By frame

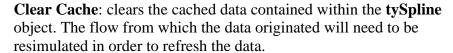
**Frame**: controls the current frame of the input spline data to display.

By speed

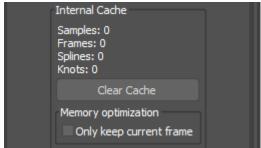
**Speed %**: the percent value that controls playback speed.

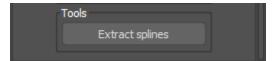
**Ref Frame**: the reference frame that the speed multiplier will be relative to.

**Internal Cache** 



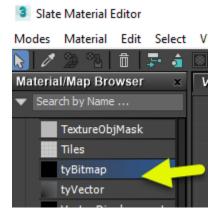
Only keep current frame: controls whether to keep data for past frames in memory, or only keep the current frame in memory. Turning this on will conserve RAM but will prevent playback of prior frames without refreshing the entire simulation. This setting is useful when rendering frames in a sequential manner on a machine with limited RAM.





# **Texmaps**

They can be accessed from the material editor



# tyVector Texmap

The **tyVector** texmap converts UVW values of sampled geometry into RGBA values whose components are the normalized lengths of individual vector axes or the vector length itself.

### TIP:

Example usage: if particles have their velocity assigned to a mapping channel, you can convert that velocity into a value between black and white using the **tyVector** texmap, depending on the length of the velocity vector for any given particle. The **tyVector** texmap can then be used as a mask for any other mixable texmap in 3ds Max, allowing you to transform particle velocities into blended color ranges...where the faster a particle travels, the more its color will change.

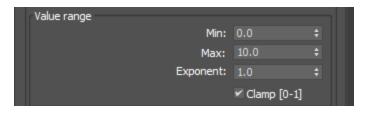


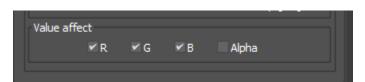
### TIP:

If the max value is greater than the min value, the interpolation will be inverted.

#### Value source

**X/Y/Z/Length**: the source of the individual float value derived from the sampled UVW value. Either the absolute value of a particular axis, or the length of the UVW vector itself.





#### Value ranges

Min/Max: the source float value will be normalized within this range, so that any value between these two values will be converted into a value between 0 and 1

**Clamp**: if enabled, normalized source values will not be clamped to 0 and 1, thus returned values may be outside of that range.

Value affect

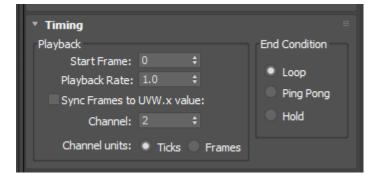
**R/G/B/A**: controls which components of the resulting RGBA value (the output of the texmap) will be set to the normalized source value. Unselected R/G/B components will be given a default value of 0. An unselected Alpha component will be given a default value of 1.

# tyBitmap Texmap

The **tyBitmap** texmap is a modified version of a default Bitmap texmap. The only difference is in the "Playback" settings of the "Timing" rollout, which contains new controls to sync bitmap sequences to user-defined particle UVW values. This allows you to control the start frame offset and playback rate on individual particles, when using animated bitmaps in your particle material.

### NOTE:

The new "Playback" controls in a **tyBitmap** will work with any UVWs on any surface, not just the particles of a **tyFlow** object. So you can bake particles to a 3rd party geometry format of your choice, and as long as the UVW coordinates are saved, the **tyBitmap** animation synchronization will still work the same.



Playback

**Start Frame**: the starting offset frame of the sequence.

**Playback Rate**: the rate at which frames will sequentially advance over time.

**Sync Frames to UVW.x value**: controls whether frame time is synchronized to the current time of the timeline, or the UVW.x value of the sampled point on the geometry's surface.

### INFO:

A UVW.x value of 0 will tell the shader to sample from frame 0 of the sequence at that particular point on the surface. A UVW.x value of 15 (if channel units is set to frames) will tell the shader to sample from frame 15 of the sequence at that particular point on the surface. Thus, by saving each particle's age to its local UVW map override within **tyFlow**, you can precisely control animation playback on each particle using this texmap.



**Channel**: controls which UVW channel to read from the geometry's surface.

### NOTE:

Make sure you choose a UVW channel that doesn't contain normal geometry UVWs (used to display diffuse/bump/etc maps) when baking timing information into your particles within **tyFlow**. The UVW channel used to synchronize timing with **tyBitmaps** won't contain proper UVWs required to actually sample the resulting images properly. For example, in most situations you would want to keep your unwrapped UVWs in channel 1, and your particle timing UVWs in channel 2.

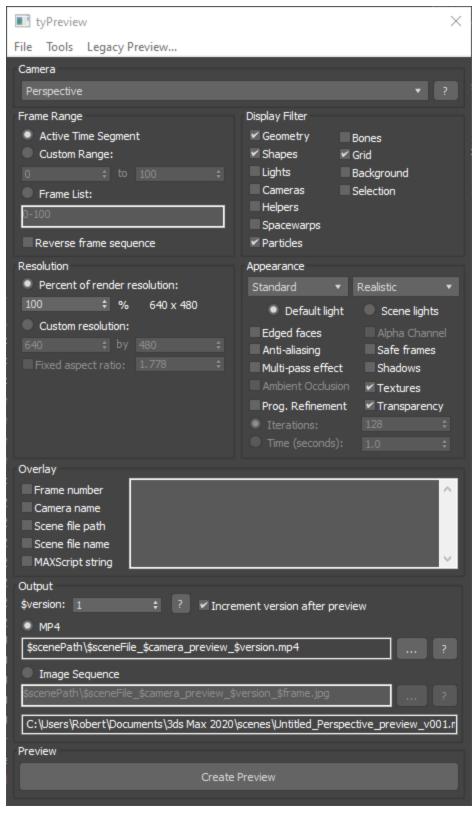


**Ticks/Frames**: controls the unit of time the channel data is saved in.

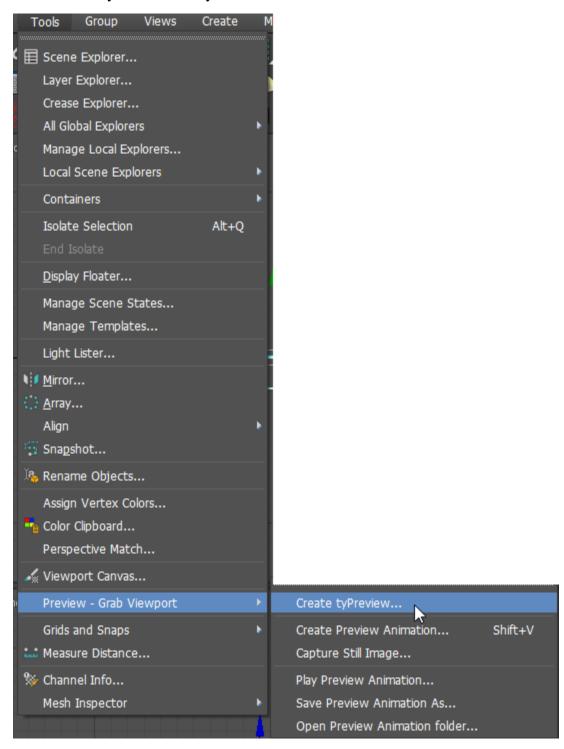
# **Utilities**

# tyPreview Utility

The **tyPreview** utility is a feature-rich replacement for 3ds Max's built-in animation preview tool.



## To access the tyPreview Utility:



### INFO:

The **tyPreview** utility has many features not available in the built-in animation preview tool. Its image processing core is multithreaded, all of its settings are saved locally to each camera/viewport, it has settings for precise output resolution control, pathname symbols, text overlays, a full MAXScript API, and it can export h264-encoded video directly to an mp4 file.

#### NOTE:

**tyPreview** does not yet multiplex audio in its output mp4 streams. If you need to preview with audio, you will still have to use 3ds Max's legacy preview tool.

### TIP:

When 3ds Max loads, **tyPreview** will automatically register itself to appropriate preview menus within Max. It also registers a macroscript (under category: **tyFlow**) that can be assigned to a custom menu or keyboard shortcut. The macroscript opens the **tyPreview** UI. To find **tyPreview** in the 3ds Max 2020 hotkey editor, simply type "tyPreview" into the action search box.

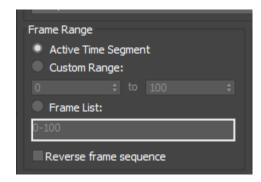


**Camera list**: this dropdown controls which camera will be previewed. Any changes made to the controls in the UI will be saved to the camera currently selected in this list. Selecting a different camera in this list will load that camera's settings into the UI.



## INFO:

All max scene object cameras (legacy camera, physical, etc) are available for preview in this list at all times. However, only currently visible built-in viewports (top, front, back, perspective, etc) are available for preview at a given time. If you wish to preview from a built-in viewport, make sure that viewport is assigned to one of the view panels first.



### Frame Range

**Active time segment**: all frames of the active time segment will be previewed.

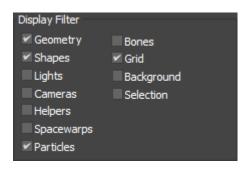
**Custom Range**: a custom range of frames will be previewed.

**Frame List**: a range of frames parsed from a user-editable textbox will be previewed.

### TIP:

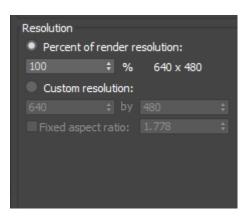
Separate groups of frames with commas, and denote ranges of frames with dashes. Ex: "10, 20, 30-50, 75-120, 125"

**Reverse frame sequence**: controls whether the preview will be made in reverse.



### **Display Filter**

**Object classes**: these checkboxes control which viewport objects and elements will be visible in the preview.



#### Resolution

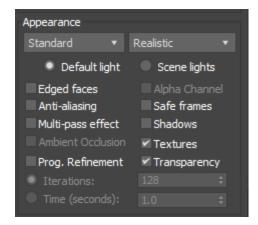
**Percent of render resolution**: the preview output resolution will be a percentage of the current render resolution.

**Custom resolution**: the preview output resolution will be an explicity-defined width and height.

**Fixed aspect ratio**: controls whether the custom resolution width/height values will be constrained to a particular aspect ratio.

### NOTE:

The 3ds Max SDK does not give low-level access to viewport resolution control, so **tyPreview** native resolution is limited to the size of a maximized viewport (on a 2K display this is roughly 1650x850 pixels). Any output resolution higher than the size of a maximized viewport will be upscaled using a bilinear filtering algorithm.



### TIP:

Use iterations mode for ensuring a certain level of progressive refinement quality is maintained per frame, regardless of how long the frame takes to render. Use time mode when you want to ensure the time taken to progressively refine frames stays under a certain threshold. Quality is not ensured in time mode, and total time is not limited in iterations mode.

### **Appearance**

**Mode**: this dropdown controls the preview rendering mode.

**Style**: this dropdown controls the preview rendering style.

**Default/scene lights**: controls whether a default light or existing scene lights will be used as a light source.

**Edged faces**: when enabled, edged faces will be visible in the preview.

**Anti-aliasing**: when enabled, 8X anti-aliasing will be enabled for the preview.

**Multi-pass effect**: when enabled, applicable multi-pass effects (DOF or motion blur) of legacy cameras will be enabled for the preview.

**Ambient Occlusion**: when enabled (with progressive refinement), ambient occlusion will be computed.

**Shadows**: when enabled, shadows will be visible.

**Prog. Refinement**: when enabled, Nitrous progressive refinement will be enabled for the preview.

**Iterations**: controls the number of progressive refinement iterations to calculate for each frame.

**Time (seconds)**: controls the amount of time to progressively refine the viewport each frame.

**Alpha channel** controls whether the preview will save alpha channel values.

**NOTE:** Alpha channel export will only work when exporting image sequences using a format that supports an alpha channel (exr, png, tif).

**Safe frames**: controls whether safe frames will be visible in the preview.

**Textures**: controls whether textures will be visible in the preview.

**Transparency**: controls whether material transparency will be visible in the preview.



**OVERLAY** 

**Frame number**: when enabled, the current frame number will be overlayed onto each frame.

**Camera name**: when enabled, the current camera name will be overlayed onto each frame.

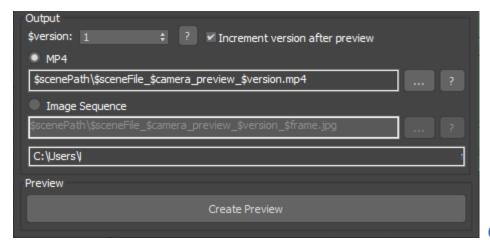
**Scene file path**: when enabled, the file's path will be overlayed onto each frame.

Scene file name: when enabled, the file's name will be overlayed onto each frame.

**MAXScript string**: when enabled, a user-defined MAXScript string will be overlayed onto each frame.

### NOTE:

The user-defined MAXScript string must evaluate to a MAXScript String object (ex: "localtime as string"). Multi-line strings are supported.



OUTPUT

### INFO:

**tyPreview** uses multiple threads to process images and save files. While viewport DIB capture is single-threaded, **tyPreview**'s gamma correction, RGB to YUV converter (mp4) and bilinear upscaling algorithms are multi-threaded. **tyPreview** also uses multiple threads to write the resulting files to disk. If you notice stuttering while exporting previews at large resolutions, that's because the multi-threaded file write function has a limiter in place which restricts the number of writer threads that can work together simultaneously. The stuttering is caused by the preview algorithm waiting for the latest batch of writes to complete before updating the time slider (in other words, stuttering is caused by disk writes happening slower than viewport updates). Stuttering is not a sign that your preview is slowing down, but instead that it's going so fast your disk cannot keep up.

**\$version**: controls the version number which will be assigned to '\$version' symbols within output path names.

**Increment version after preview**: when enabled, the version number will be incremented by 1 after each successfully completed preview. Canceled/failed previews will not trigger an increment.

**MP4**: when selected, the preview output will be saved to an h264-encoded mp4 file.

### INFO:

No mp4 encoder settings are exposed in the **tyPreview** UI. The hard-coded mp4 encoder settings are designed to achieve a reasonable bitrate while reducing artifacts in exports.

**MP4 path**: the output path where mp4 files will be saved.

**Image Sequence**: when selected, the preview output will be saved to a sequence of image files.

### INFO:

The available image-sequence filetypes are jpg, png, tif and exr. No image encoder settings are exposed in the **tyPreview** UI. All hard-coded image encoder settings are set to ensure maximum output quality, and alpha channel output (where applicable).

**Image sequence path**: the output path where images will be saved.

### TIP:

Path values can contain various symbols to denote built-in path variables. Click the "?" button beside each path box to see the list of available symbols. The textbox below the editable path inputs contains the resulting fully-parsed output path.

## **MAXScript access**

• **tyPreviewWindow**(): this function loads the **tyPreview** UI.

The **tyPreview** utility can also be executed purely through MAXScript, without having to open the UI.

• **tyPreview**(): this function executes a **tyPreview** export, using settings for the currently active viewport.

For **tyPreview** customization, the **tyPreview** MAXScript API gives you full control over all **tyPreview** settings. All of these settings can be called as arguments for the base **tyPreview()** MAXScript function. For example, instead of "**tyPreview()**", you might call "**tyPreview frameRange\_reverse:true displayFilter\_geometry:false appearance\_alpha:true**".

#### TIP:

The "tyPreviewWindow()" function can also be called with all of the same function arguments available to the "tyPreview()" function. Instead of immediately executing a preview with those arguments, it will adjust controls in the UI to match your inputs.

All of the **tyPreview** function arguments correspond to settings already described in the documentation above, so while they will be listed below, redundant descriptions for them are not included.

### tyPreview function arguments:

- **camera\_name**: (name of target camera node)
- camera node (target camera node)

### NOTE:

If camera\_node is specified, tyPreview will attempt to preview from that camera. If camera\_name is specified, tyPreview will attempt to preview from the camera with that name. If neither is specified, tyPreview will preview from the camera of the current active viewport. Export settings will be taken from whichever camera is ultimately chosen. Those settings can be overridden using the function arguments listed below.

- frameRange\_list: (ex: "10, 20, 30-50, 75-120, 125")
- frameRange\_reverse:
- displayFilter\_geometry: true|false
- displayFilter\_shapes: true|false
- displayFilter\_lights: true|false
- displayFilter\_cameras: true|false
- displayFilter\_helpers: true|false
- displayFilter\_spacewarps: true|false
- displayFilter\_particles: true|false
- displayFilter bones: true|false
- displayFilter\_grid: true|false
- displayFilter\_background: true|false
- resolution width: integer
- resolution\_height: integer
- appearance\_mode: integer (0-based index of modes listed in tyPreview UI)
- appearance\_style: integer (0-based index of styles listed in tyPreview UI)
- appearance\_lights: integer (0 = default; 1 = scene)
- appearance\_edgedFaces: true|false
- appearance\_alpha: true|false
- appearance\_antiAliasing: true|false
- appearance\_safeFrames: true|false
- appearance\_multipass: true|false
- appearance\_textures: true|false
- appearance\_transparency: true|false
- appearance\_ao: true|false
- appearance\_shadows: true|false

- appearance\_progressiveRefinement: true|false
- **appearance\_progressiveRefinement\_type**: integer (0 = iterations; 1 = time)
- appearance\_progressiveRefinement\_iterations: integer
- appearance\_progressiveRefinement\_time: float
- **overlay\_frame**: true|false
- overlay\_camera: true|false
- overlay\_sceneFile: true|false
- overlay\_scenePath: true|false
- **overlay\_maxscript**: true|false
- overlay\_maxscript\_script: string
- output\_version: integer
- **output\_type**: integer (0 = mp4; 1 = images)
- output\_framerate: float

### NOTE:

output\_type:1 must be specified for image sequence export.

• output\_filename: string